# On perturbation around closed exclusion processes

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## Abstract

We derive the formula for the stationary states of particle-number conserving exclusion processes infinitesimally perturbed by inhomogeneous adsorption and desorption. The formula not only proves but also generalises the conjecture proposed in [Phys. Rev. E 97, 032135] to account for inhomogeneous adsorption and desorption. As an application of the formula, we draw part of the phase diagrams of the open asymmetric simple exclusion process with and without Langmuir kinetics, correctly reproducing known results.

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# 15 1 Introduction

Among the famous solvable models of driven diffusive systems is the asymmetric simple exclusion process (ASEP). Aside from being solvable deep in the non-equilibrium regime, the
model is interesting for its connections to various ideas in statistical physics such as boundaryinduced phase transitions [1], the KPZ universality class [2], and random matrix theory [3]. It

has also attracted attention for its wide applicability to various phenomena in physics, biology
 or society [4–6].

ASEP describes particles on a one-dimensional lattice that hop asymmetrically (e.g., more 22 frequently to the right than left). It is a continuous-time Markov process and as such it is in-23 teresting to study where the probability distribution settles after a long time, *i.e.*, its stationary 24 distribution (state). This is equivalent to studying the eigenvector(s) of the Markov matrix 25 with zero eigenvalue, which is possible for ASEP because of its  $U_q(sl(2))$  symmetry [7–9]. 26 However, remained elusive is the stationary distribution (as well as other properties) for ex-27 clusion processes without integrability, even though their practical usefulness and applicability 28 are not impaired by the lack thereof. 29

One example of exclusion processes without integrability is the ASEP combined with Langmuir kinetics (ASEP-LK), where, in addition to asymmetric hopping, a particle can attach to or detach from the lattice at homogeneous rates. Even though the system has the  $U_q(sl(2))$ symmetry on a periodic chain, it is indeed broken on a finite-length chain with boundary conditions. The model is proposed to describe unidirectional motion of motor proteins [10], so it is a good example of non-integrable exclusion processes with interesting applications.

One possible way to analytically study the properties of non-integrable exclusion processes is to take the thermodynamic limit. One can for example determine the phase diagram of the system by solving the fluid equations obtained by taking the coarse-grained, continuum limit. However, the strategy usually involves using the mean-field approximation, which may or may not be justified from first principles. It would be desirable if we have another theoretical method with a clear regime of validity to compare with experiments or computer simulations. This could theoretically justify the mean-field approximations as well.

There is one such method which seemingly has been mostly overlooked – the perturbation theory. The ASEP-LK is a prime example: The stationary states of the periodic/closed/open ASEP have been obtained exactly, and so one can in principle obtain those of the ASEP-LK perturbatively when the ad/desorption rates are small. For example, [11] conjectured a formula for the stationary states of the closed ASEP with infinitesimally small Langmuir kinetics. This formula is yet to be proven despite having a simple and inviting form, however.

The goal of this paper is to set up such perturbation in generic situations. We consider 49 perturbing a particle-number conserving hopping process with inhomogeneous ad/desorption 50 and derive a formula for the stationary state at leading order. (The leading order result is 51 meaningful because this is a degenerate perturbation theory.) Our formula potentially has 52 various interesting applications. For one thing, the above-mentioned conjecture is its imme-53 diate consequence since closed ASEP is clearly a particle-number conserving process. We can 54 also apply the formula to draw perturbative part of the phase diagrams of the open ASEP 55 with/without Langmuir kinetics. We do so by interpreting the open boundary condition as a 56 special case of the inhomogeneous ad/desorption acting only on boundary sites. The results, 57 as we will see later, reproduce the results obtained in [12] but without relying on the mean-58 field approximation or any other unjustified approximations at all. Therefore we are going to 59 have a clear regime of validity for our theoretical formula, even though the price we pay is the 60 restriction to the perturbative regime. 61

The rest of the paper is organised as follows. We first briefly review Markov processes, in particular closed ASEP with/without infinitesimal Langmuir kinetics in Section 2. We then go on to construct the stationary states of generic particle-number conserving Markov processes infinitesimally perturbed by inhomogeneous ad/desorption in Section 3. This will, as a special case, prove the conjecture given in [11]. We will provide other applications of the formula by deriving the phase diagram of the open ASEP with and without Langmuir kinetics in Section 4. We conclude in Section 5 with discussions and future directions.

## <sup>69</sup> 2 Driven diffusive systems with ad/desorption

#### 70 2.1 Continuous-time Markov process

<sup>71</sup> Let us consider a continuous-time Markov process describing particles hopping on *L* lattice <sup>72</sup> sites (of arbitrary shapes and dimensions). We are interested in the time evolution of the <sup>73</sup> probability associated with a given configuration. This is given by a collection of differential <sup>74</sup> equations, conveniently written in matrix form using the Markov matrix *M*,

$$\frac{d}{dt} \left| P \right\rangle = M \left| P \right\rangle, \tag{2.1}$$

where  $|P\rangle$  is a vector collecting probabilities of realising given configurations. In other words, by writing the configuration of particles as  $(\tau_1, ..., \tau_L)$  where  $\tau_i = 1$  ( $\tau_i = 0$ ) means that a

particle is (not) present at site *i*, and its realisation probability as  $p(\tau_1, \ldots, \tau_L)$ , we package

78 the distribution into a state

$$|P\rangle = \sum_{\tau_1,\dots,\tau_L} p(\tau_1,\dots,\tau_L) |\tau_1,\dots,\tau_L\rangle, \qquad (2.2)$$

<sup>79</sup> and this vector evolves according to the differential equation above.

For later convenience we denote the total Hilbert space as V, which is a tensor product of the Hilbert space  $V_i$  on site i from i = 1 to L,

$$V = \bigotimes_{i=1}^{L} V_i. \tag{2.3}$$

<sup>82</sup> It can also be decomposed into a direct sum of fixed particle number subspaces  $W_N$  (where N

<sup>83</sup> indicates the number of particles in the system), so that

$$V = \bigoplus_{N=0}^{L} W_N. \tag{2.4}$$

Given such an evolution equation, one interest lies in finding where the probability distribution settles after a long time. This is given by the eigenvector of M with eigenvalue zero. The number of such eigenvectors are expected to match that of the superselection sectors of M.

#### <sup>88</sup> 2.2 Closed exclusion process with ad/desorption

<sup>89</sup> Our interest in this paper lies in the system *M* such that  $M = M_0 + \epsilon H$  where  $\epsilon \ll 1$  is <sup>90</sup> the perturbation parameter<sup>1</sup>. We require that  $M_0$  conserves the particle number (*i.e.*, U(1)<sup>91</sup> symmetric) while  $\epsilon H$  breaks it *via* ad/desorption. Concretely, we have

$$M \equiv M_0 + \epsilon H, \quad M_0 \Big|_{W_N} : W_N \to W_N$$
  
$$H = \sum_{i=1}^L h_i, \quad h_i \equiv \begin{pmatrix} -\alpha_i & \beta_i \\ \alpha_i & -\beta_i \end{pmatrix}_{V_i}.$$
 (2.5)

where  $h_i$  only acts on *i*-sites, with  $\epsilon \alpha_i$  and  $\epsilon \beta_i$  representing adsorption and desorption rates at site *i*, respectively.

Note that, before perturbation, the state space breaks up into L + 1 superselection sectors  $W_N$ , and we have a stationary state for each of them. We denote the *N*-particle stationary state

<sup>&</sup>lt;sup>1</sup>We will hereafter identify the Markov matrix as the corresponding system itself.

<sup>96</sup> as  $|S_N\rangle$  hereafter. In addition, since the perturbation breaks the particle-number symmetry, <sup>97</sup> there are no more superselection sectors for *M*. It therefore has only one stationary state, <sup>98</sup> which we denote as  $|\tilde{S}\rangle$ .

The goal of this paper is to construct  $|\tilde{S}\rangle$  in terms of  $|S_N\rangle$  at leading order in  $\epsilon$ . It is by now clear that this is the zeroth order degenerate perturbation theory. We need to find the right basis on which the higher-order perturbation theory is run. We will study this in Section 3.

Example: ASEP with Langmuir kinetics Before moving on, we present an example of such
 systems, known as the closed ASEP perturbed by Langmuir kinetics (ASEP-LK). The Markov
 matrix of closed ASEP-LK is given by the following,

$$M = M_0 + \epsilon H$$

$$M_0 = \sum_{i=1}^{L-1} M_{i,i+1}, \quad M_{i,i+1} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & -q & 1 & 0 \\ 0 & q & -1 & 0 \\ 0 & 0 & 0 & 0 \end{pmatrix}_{V_i \otimes V_{i+1}}$$

$$H = \sum_{i=1}^{L} h_i, \quad h_i \equiv \begin{pmatrix} -\alpha & 1 \\ \alpha & -1 \end{pmatrix}_{V_i}.$$
(2.6)

where  $M_0$  describes the closed ASEP and  $\epsilon H$  the Langmuir kinetics.  $M_{i,i+1}$  acts as an identity outside  $V_i \otimes V_{i+1}$  and the bases of  $M_{i,i+1}$  inside  $V_i \otimes V_{i+1}$  are given by, from top to bottom columns or left to right rows,  $|0,0\rangle$ ,  $|0,1\rangle$ ,  $|1,0\rangle$ , and  $|1,1\rangle$ . The perturbation  $\epsilon H$  describes a particle homogeneously detaching from the lattice at rate  $\omega_d \equiv \epsilon$  while attaching at rate  $\omega_a \equiv \epsilon \alpha$ .

The closed ASEP,  $M_0$ , trivially conserves the particle number and so has superselection sectors labelled by it. There are therefore L+1 stationary states in  $M_0$ , which can be computed by using the Bethe ansatz as [13]

$$|S_N\rangle = \begin{bmatrix} L\\N \end{bmatrix}_q^{-1} \sum_{(n)_N} q^{\sum_{j=1}^N (L-j+1-n_j)} |(n)_N\rangle, \qquad (2.7)$$

where  $|S_N\rangle$  denotes the *N*-particle stationary state. Here  $\begin{bmatrix} L \\ N \end{bmatrix}_q$  is the *q*-binomial, defined by

$$\begin{bmatrix} L \\ N \end{bmatrix}_{q} \equiv \frac{(q;q)_{L}}{(q;q)_{N}(q;q)_{L-N}}, \quad (a;q)_{n} \equiv \prod_{i=1}^{n} (1-aq^{i-1}), \quad (2.8)$$

and  $(n)_N$  is an ordered collection of N sites on which the particles are present. We also used a shorthand notation  $|(n)_N\rangle$  to refer to the basis corresponding to such a configuration. The overall normalisation is because the sum of probabilities must be one.

Once we perturb the system by  $\epsilon H$ , the particle-number conservation is lost and there is only one stationary state,  $|\tilde{S}\rangle$ . Because the integrability is (mostly likely) lost due to the perturbation, it is considered difficult to derive the stationary state for this model. It was however conjectured in [11] that in the  $\epsilon \equiv \omega_d \rightarrow 0$  limit (while fixing  $\alpha$ )  $|\tilde{S}\rangle$  is given by

$$|\tilde{S}\rangle = \frac{1}{(1+\alpha)^L} \sum_{N=0}^{L} {L \choose N} \alpha^N |S_N\rangle + O(\epsilon)$$
(2.9)

<sup>121</sup> We will prove this conjecture in the next section as a corollary to the main result.

## <sup>122</sup> **3** Construction of the stationary state

- <sup>123</sup> We are going to prove the following theorem.
- 124 **Theorem 1.** For a class of continuous-time Markov processes  $M = M_0 + \epsilon H$  defined in (2.5), the
- stationary state of M can be written in terms of the N-particle stationary states of  $M_0$ ,  $|S_N\rangle$ , as

$$|\tilde{S}\rangle \equiv \frac{1}{\sum_{N=0}^{L} p_N} \sum_{N=0}^{L} p_N |S_N\rangle + O(\epsilon), \quad p_N = \prod_{i=1}^{N} \left(\frac{A_L - A_{i-1}}{B_i}\right).$$
(3.1)

where  $A_N$  and  $B_N$  are given by

$$A_N \equiv \sum_{(n)_N} q[(n)_N] \sum_{n \in (n)_N} \alpha_n, \qquad (3.2)$$

$$B_N \equiv \sum_{(n)_N} q[(n)_N] \sum_{n \in (n)_N} \beta_n, \qquad (3.3)$$

127 using

$$|S_N\rangle \equiv \sum_{(n)_N} q[(n)_N] |(n)_N\rangle, \quad \sum_{(n)_N} q[(n)_N] = 1$$
 (3.4)

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<sup>129</sup> Incidentally, we have

$$A_0 = 0, \quad A_L = \sum_{i=1}^{L} \alpha_i, \quad B_0 = 0, \quad B_L = \sum_{i=1}^{L} \beta_i$$
 (3.5)

Before attempting to prove this theorem, we have one remark.

**Corollary 1.** This, as a corollary, immediately proves the conjecture (2.9) given in [11].

Proof of Corollary 1. Because  $A_N = N\alpha$  and  $B_N = N$  in the current case, we immediately have  $p_N = {L \choose N} \alpha^N$ . We therefore have  $\sum_{N=0}^{L} p_N = (1 + \alpha)^L$ . This concludes the proof of the conjecture (2.9).

Now we move on to proving Theorem 1, but prior to this let us set up some notations which will be useful later. We denote  $K_0$  as the subspace spanned by all the stationary states of  $M_0$ , while  $K_1$  as the subspace spanned by all other eigenvectors. Because  $M_0$  is non-normal,  $K_0$ and  $K_1$  are not orthogonal to each other.

Let us also present a general argument to understand the strategy of the proof. Notice that we are trying to find the eigenvector of a matrix in perturbation theory, starting from degenerate vacua. In order to do this, we need to find a vector in  $K_0$  which is taken to  $K_1$  upon acting with  $\epsilon H$ . (For example see Appendix A.4 of [14].) This will single out a linear combination of the stationary states of  $M_0$ , on which higher-order perturbations can be studied. Put differently, we need to find  $|S\rangle$  such that

$$|S\rangle \in K_0 \quad \text{and} \quad H|S\rangle \in K_1,$$
 (3.6)

where  $|S\rangle$  denotes the  $O(\epsilon^0)$  part of  $|\tilde{S}\rangle$ .

It might seem as if one needs to know all the eigenvectors of  $M_0$  in order to impose such conditions. However, this is too pessimistic. The space  $K_1$  can be characterised by the fact that its inner product with the left eigenvector of  $M_0$  with vanishing eigenvalue is zero. In

other words, if we write  $|L_N\rangle$  as the *N*-particle eigenvector of  $M_0^T$  (the transpose of  $M_0$ ) with 149 vanishing eigenvalue, 150

$$M_0^T \left| L_N \right\rangle = 0, \tag{3.7}$$

we have that 151

$$\langle L_N | \psi \rangle = 0 \quad \text{if} \quad | \psi \rangle \in K_1.$$
 (3.8)

In addition, the form of  $|L_N\rangle$  is immediate because  $M_0$  is a Markov matrix, 152

$$|L_N\rangle = \sum_{(n_i)_N} |(n_i)_N\rangle.$$
(3.9)

This hinges on the fact that the sum of probabilities is constant in time and hence the sum of 153 columns in a Markov matrix is zero (in each superselection sector, if any).<sup>2</sup> The normalisation 154 is immaterial so we chose an arbitrary one. 155

Summarising the discussions above, we now need to find  $|S\rangle \in K_0$  such that  $\langle L_N | H | S \rangle = 0$ 156

for any N. We parameterise  $|S\rangle$  for convenience as 157

$$|S\rangle \equiv \frac{1}{\sum_{N=0}^{L} p_N} \sum_{N=0}^{L} p_N |S_N\rangle, \qquad (3.10)$$

where we can set  $p_0 = 1$ . We also parameterise  $|S_N\rangle$  as 158

$$|S_N\rangle \equiv \sum_{(n)_N} q[(n)_N] |(n)_N\rangle.$$
(3.11)

We demand that they are properly normalised, in other words that the sum of probabilities 159 becomes one,  $\sum_{(n)_N} q[(n)_N] = 1$ . Let us prove Theorem 1 now. 160

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*Proof of Theorem 1.* First of all,  $H|S_i\rangle$  does not overlap with  $|L_N\rangle$  unless i = N-1, N, or N+1162 because *H* only takes *i*-particle states to *i*- or  $(i \pm 1)$ -particle states. Therefore the conditions 163  $\langle L_N | H | S \rangle = 0$  reduce to a set of recursion relations, 164

$$p_{N-1} \langle L_N | H | S_{N-1} \rangle + p_N \langle L_N | H | S_N \rangle + p_{N+1} \langle L_N | H | S_{N+1} \rangle = 0, \qquad (3.12)$$

where we set  $p_{-1} = p_{L+1} = 0$  for consistency. 165

Let us now compute  $\langle L_N | H | S_i \rangle$  for i = N - 1, i = N, and i = N + 1. Because we only 166 need to compute the overlap with  $|L_N\rangle$ , we will only compute the projection of  $H|S_i\rangle$  to  $W_N$ . 167 Starting from i = N - 1, we have 168

$$H|S_{N-1}\rangle \bigg|_{W_N} = \sum_{(n)_{N-1}} q[(n)_{N-1}] \sum_{n \notin (n)_{N-1}} \alpha_n |(n)_{N-1} \cup n\rangle, \qquad (3.13)$$

where  $|(n)_{N-1} \cup n|$  means adding a particle on site *n* to the (N-1)-particle state  $|(n)_{N-1}\rangle$ . 169 We then have 170

$$\langle L_N | H | S_{N-1} \rangle = \sum_{(n)_{N-1}} q[(n)_{N-1}] \sum_{n \notin (n)_{N-1}} \alpha_n$$
 (3.14)

$$= \sum_{(n)_{N-1}} q[(n)_{N-1}] \left( \sum_{n=1}^{L} \alpha_n - \sum_{n \in (n)_{N-1}} \alpha_n \right) = A_L - A_{N-1}$$
(3.15)

<sup>&</sup>lt;sup>2</sup>The form of  $|L_N\rangle$  suggests that the overlap  $\langle L_N | \psi \rangle$  is the sum of probabilities of realising N-particle states in  $|\psi\rangle$ . We thank Yuki Ishiguro and Jun Sato for discussions on this point. See also their paper [15] whose submission was coordinated with ours.

Let us continue to i = N. The *N*-particle subspace component in  $H |S_N\rangle$  is given by

$$H|S_N\rangle\Big|_{W_N} = -\sum_{(n)_N} q[(n)_N] \sum_{n \notin (n)_N} \alpha_n |(n)_N\rangle - \sum_{(n)_N} q[(n)_N] \sum_{n \in (n)_N} \beta_n |(n)_N\rangle$$
(3.16)

<sup>172</sup> The overlap with  $|L_N\rangle$  is hence given by

$$\langle L_N | H | S_N \rangle = -(A_L - A_N + B_N) \tag{3.17}$$

Finally we study the case where i = N+1. The *N*-particle subspace component in  $H |S_{N+1}\rangle$ is given by

$$H|S_{N+1}\rangle \bigg|_{W_N} = \sum_{(n)_{N+1}} q[(n)_{N+1}] \sum_{n \in (n)_{N+1}} \beta_n |(n)_{N+1} \setminus n\rangle$$
(3.18)

where  $|(n_i) \setminus n\rangle$  means removing a particle on site n from the (N + 1)-particle state  $|(n)_{N+1}\rangle$ . The overlap with  $|L_N\rangle$  is hence given by

$$\langle L_N | H | S_N \rangle = B_{N+1} \tag{3.19}$$

177 The recursion relation (3.12) therefore becomes

$$(A_L - A_{N-1})p_{N-1} - B_N p_N = (A_L - A_N)p_N - B_{N+1}p_{N+1}.$$
(3.20)

Since we have  $(A_L - A_{N-1})p_{N-1} - B_N p_N|_{N=0} = 0$ , we can derive a simplified recursion relation, tion,

$$p_N = \frac{A_L - A_{N-1}}{B_N} p_{N-1}.$$
(3.21)

By solving this recursion relation, we conclude that the stationary state of the system Mbecomes

$$|\tilde{S}\rangle \equiv \frac{1}{\sum_{N=0}^{L} p_{N}} \sum_{N=0}^{L} p_{N} |S_{N}\rangle + O(\epsilon), \quad p_{N} = \prod_{i=1}^{N} \left(\frac{A_{L} - A_{i-1}}{B_{i}}\right).$$
(3.22)

<sup>182</sup> In other words we have successfully proven Theorem 1.

# <sup>183</sup> 4 Phase diagram of the open ASEP(-LK)

It is interesting to apply our formula to derive the phase diagram of the open ASEP with/without
Langmuir kinetics in terms of perturbation theory. This can be done by considering the open
boundary condition as a particular case of the inhomogeneous ad/desorption. More conretely,
the open ASEP-LK is defined by the following Markov matrix

$$M \equiv M_0 + \tilde{H}, \quad \tilde{H} = \sum_{i=1}^L \tilde{h}_i, \quad \tilde{h}_i \equiv \begin{pmatrix} -\boldsymbol{\omega}_i^{[a]} & \boldsymbol{\omega}_i^{[d]} \\ \boldsymbol{\omega}_i^{[a]} & -\boldsymbol{\omega}_i^{[d]} \end{pmatrix}_{V_i}, \quad (4.1)$$

where  $M_0$  is the Markov matrix of the closed ASEP, while we demand  $\omega_2^{[a]} = \omega_3^{[a]} = \cdots = \omega_{L-1}^{[a]}$  and  $\omega_2^{[d]} = \omega_3^{[d]} = \cdots = \omega_{L-1}^{[d]}$ . Note that the system becomes the open ASEP without Langmuir kinetics when  $\omega_i^{[a]} = \omega_i^{[d]} = 0$  for i = 2, ..., L - 1. When  $\omega_i^{[a]}$  and  $\omega_i^{[d]}$  are small, the system is amenable to perturbation theory and our formula (3.22) is applicable. We
 therefore set

$$\omega_{1}^{[a]} = \epsilon \alpha, \quad \omega_{L}^{[d]} = \epsilon \beta$$
  

$$\omega_{1}^{[d]} = \epsilon \gamma, \quad \omega_{L}^{[a]} = \epsilon \delta$$
  

$$\omega_{i}^{[a]} = \epsilon a, \quad \omega_{i}^{[d]} = \epsilon b \quad \text{for } i = 2, \dots, L-1$$
(4.2)

and compute the stationary state of the open ASEP-LK at leading order in  $\epsilon \ll 1$ . In other words, we have, in the language of (2.5),

$$a_1 = \alpha, \quad a_2 = \dots = a_{L-1} = a, \quad a_L = \delta$$
  

$$\beta_1 = \gamma, \quad \beta_2 = \dots = \beta_{L-1} = b, \quad \beta_L = \beta$$
(4.3)

<sup>195</sup> For later convenience, we will denote the *N*-particle stationary state of the closed ASEP as

$$|S_N\rangle \equiv \sum_{(n)_N} q_L[(n)_N] |(n)_N\rangle, \quad q_L[(n)_N] = {L \brack N}_q^{-1} q^{\sum_{j=1}^N (L-j+1-n_j)}$$
(4.4)

emphasising that the number of sites is *L*. We will also denote  $q_L[(n)_N | \tau_1 = 0, 1, \tau_L = 0, 1]$ to restrict  $(n)_N$  to obey particles at site 1 and *L* being present/absent. Equivalently, we can set  $q_L[(n)_N | \tau_1 = 0, 1, \tau_L = 0, 1] = 0$  if  $(n)_N$  is not consistent with  $\tau_1 = 0, 1$  or  $\tau_L = 0, 1$ .

Let us now compute  $A_i$  and  $B_i$ . We hereafter restrict our attention to  $A_i$  only since  $B_i$  can be obtained from  $A_i$  by swapping a with  $\gamma$ ,  $\delta$  with  $\beta$ , and a with b. We have

$$A_N = \sum_{\tau_1, \tau_L = 0, 1} A_N^{\tau_1, \tau_L}$$
(4.5)

where (for example)  $A_N^{0,1}$  means that the sum over  $(n)_N$  in the definition of  $A_N$  is restricted to its subset in which  $\tau_1 = 0$  (absent) and  $\tau_L = 1$  (present). More concretely, they are defined as

$$A_N^{\tau_1,\tau_L} \equiv A_N \equiv \sum_{(n)_N} q_L[(n)_N | \tau_1, \tau_L] \sum_{n \in (n)_N} \alpha_n$$
(4.6)

This will not become too complicated as  $q_L[(n)_N | \tau_1, \tau_L]$  can be related to  $q_{L-2}[(n)_N]$ ,  $q_{L-2}[(n)_{N-1}]$ , etc. For example,

$$q_{L}[(n)_{N}|\tau_{1}=0,\tau_{L}=0]=q^{2N-N}q_{L-2}[(n)_{N}], \qquad (4.7)$$

where 2*N* and -N in the exponent comes from the shifting of *L* to L-2 and of  $n_j$  to  $n_j - 1$ , respectively. Similar arguments lead to

$$q_{L}[(n)_{N} | \tau_{1} = 0, \tau_{L} = 0] = q^{N} \times q_{L-2}[(n)_{N}]$$

$$q_{L}[(n)_{N} | \tau_{1} = 0, \tau_{L} = 1] = q^{0} \times q_{L-2}[(n)_{N-1}],$$

$$q_{L}[(n)_{N} | \tau_{1} = 1, \tau_{L} = 0] = q^{L-1} \times q_{L-2}[(n)_{N-1}],$$

$$q_{L}[(n)_{N} | \tau_{1} = 1, \tau_{L} = 1] = q^{L-N} \times q_{L-2}[(n)_{N-2}],$$
(4.8)

## 208 We therefore have

$$A_{N}^{0,0} = \frac{{\binom{L-2}{N}}_{q}}{{\binom{L}{N}}_{q}} q^{N} \times aN$$

$$A_{N}^{0,1} = \frac{{\binom{L-2}{N-1}}_{q}}{{\binom{L}{N}}_{q}} q^{0} \times (a(N-1)+\delta)$$

$$A_{N}^{1,0} = \frac{{\binom{L-2}{N-1}}_{q}}{{\binom{L}{N}}_{q}} q^{L-1} \times (a+a(N-1))$$

$$A_{N}^{1,1} = \frac{{\binom{L-2}{N-2}}_{q}}{{\binom{L}{N}}_{q}} q^{L-N} \times (a+a(N-2)+\delta).$$
(4.9)

209 and likewise

$$B_{N}^{0,0} = \frac{{\binom{L-2}{N}}_{q}}{{\binom{L}{N}}_{q}} q^{N} \times bN$$

$$B_{N}^{0,1} = \frac{{\binom{L-2}{N-1}}_{q}}{{\binom{L}{N}}_{q}} q^{0} \times (b(N-1) + \beta)$$

$$B_{N}^{1,0} = \frac{{\binom{L-2}{N-1}}_{q}}{{\binom{L}{N}}_{q}} q^{L-1} \times (\gamma + b(N-1))$$

$$B_{N}^{1,1} = \frac{{\binom{L-2}{N-2}}_{q}}{{\binom{L}{N}}_{q}} q^{L-N} \times (\gamma + b(N-2) + \beta).$$
(4.10)

# 210 4.1 Phase diagram of the open ASEP

<sup>211</sup> We are now ready to compute the stationary state of the open ASEP by setting  $a = b = \gamma =$ <sup>212</sup>  $\delta = 0$ . From the above computations, we have

$$A_i = \alpha q^{L-N} \times \frac{1-q^N}{1-q^L}, \quad B_i = \beta \times \frac{1-q^N}{1-q^L},$$
 (4.11)

<sup>213</sup> which leads to

$$p_N \equiv \prod_{i=1}^N \left( \frac{A_L - A_{i-1}}{B_i} \right) = \left( \frac{\alpha}{\beta} \right)^N \times \begin{bmatrix} L \\ N \end{bmatrix}_q$$
(4.12)

Therefore the stationary state  $|\tilde{S}\rangle$  of the open ASEP becomes, at leading order in  $O(\epsilon)$ ,

$$|\tilde{S}\rangle \equiv \frac{1}{\sum_{N=0}^{L} (\alpha/\beta)^{N} \times {L \brack N}_{q}} \sum_{N=0}^{L} \left(\frac{\alpha}{\beta}\right)^{N} \times {L \brack N}_{q} |S_{N}\rangle + O(\epsilon)$$
(4.13)

$$= \frac{1}{\sum_{N=0}^{L} (\alpha/\beta)^N \times {L \brack N}_q} \sum_{N=0}^{L} \left(\frac{\alpha}{\beta}\right)^N \sum_{(n)_N} q^{\sum_{j=1}^{N} (L-j+1-n_j)} |(n)_N\rangle + O(\epsilon), \qquad (4.14)$$

from which all relevant physical quantities (particle number density, *n*-point functions, etc.) 215 can be computed. Incidentally, the normalisation constant can be written more compactly as 216

$$\sum_{N=0}^{L} \left(\frac{\alpha}{\beta}\right)^{N} \times \begin{bmatrix} L\\ N \end{bmatrix}_{q} = {}_{2}\phi_{0} \begin{bmatrix} q^{-N}, 0\\ \emptyset \end{bmatrix}; q, \frac{\alpha}{\beta} \times q^{N} \end{bmatrix}$$
(4.15)

where  $_{r}\phi_{s}$  is the *q*-hypergeometric function, defined as 217

$${}_{r}\phi_{s}\begin{bmatrix}a_{1},a_{2},\ldots,a_{r}\\b_{1},b_{2},\ldots,b_{s};q,z\end{bmatrix} = \sum_{n=0}^{\infty} \frac{(a_{1},a_{2},\ldots,a_{r};q)_{n}}{(b_{1},b_{2},\ldots,b_{s},q;q)_{n}} \left((-1)^{n}q^{\binom{n}{2}}\right)^{s+1-r} z^{n}, \qquad (4.16)$$

218

in which  $(a_1, a_2, ..., a_r; q)_n \equiv \prod_{i=1}^r (a_i; q)_n$ . Let us now detect the phase transition in the open ASEP by computing the particle number 219 density, or equivalently, the one point function  $\langle \tau_i \rangle$ . For the sake of simpler analytic computa-220 tions, we hereafter restrict our attention to q = 0. This makes thing particularly easy because 221 the particle number density  $\langle \tau_i \rangle_N$  of  $|S_N\rangle$  is given by the step function, 222

$$\langle \tau_i \rangle_N = \begin{cases} 1 & i \ge L - N + 1 \\ 0 & i \le L - N \end{cases}.$$

$$(4.17)$$

The number density  $\langle \tau_i \rangle$  of  $|\tilde{S}\rangle$  is then given by (at leading order in  $\epsilon$ ) 223

$$\langle \tau_i \rangle = \frac{\sum_{N=L-i+1}^{L} (\alpha/\beta)^N}{\sum_{N=0}^{L} (\alpha/\beta)^N} = \frac{(\alpha/\beta)^{L+1-i} - (\alpha/\beta)^{L+1}}{1 - (\alpha/\beta)^{L+1}}$$
(4.18)

where we used  $\lim_{q\to 0} {L \brack N}_q = 1$ . We plot  $\langle \tau_i \rangle$  for some values of  $\alpha/\beta$  in Figure 1. 224

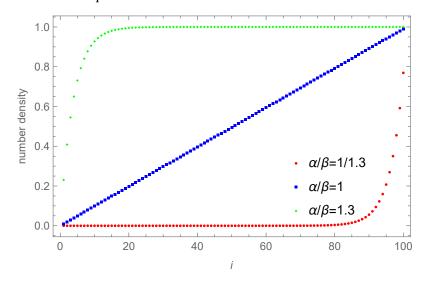


Figure 1: Plot of the particle number densities of the open ASEP (at q = 0) as functions of lattice sites *i*. We take the number of lattice sites to be L = 100. For *L* as large as 100, we already see three distinct phases –  $\alpha/\beta$  < 1 corresponds to the low-density phase,  $\alpha/\beta = 1$ , the coexistence phase, and  $\alpha/\beta > 1$ , the high-density phase.

Let us take the thermodynamic limit  $L \to \infty$ . It is immediate to see that the behaviour of 226  $\langle \tau_i \rangle$  are completely different for three cases,  $\alpha/\beta \leq 1$ . For  $\alpha/\beta < 1$ , we have

$$\langle \tau_i \rangle = \begin{cases} 0 & \text{for } L - i \gg L^0 \\ \left(\frac{\alpha}{\beta}\right)^{L+1-i} & \text{for } L - i = O(L^0) \end{cases}$$
(4.19)

227 for  $\alpha/\beta = 1$ ,

$$\langle \tau_i \rangle = \frac{i}{L+1},\tag{4.20}$$

<sup>228</sup> and for  $\alpha/\beta > 1$ ,

$$\langle \tau_i \rangle = \begin{cases} 1 - \left(\frac{\alpha}{\beta}\right)^{-i} & \text{for } i = O(L^0) \\ 1 & \text{for } i \gg O(L^0) \end{cases}.$$
(4.21)

<sup>229</sup> Corresponding to the number density in the bulk region of the open chain, we call the phase <sup>230</sup> realised for  $\alpha/\beta < 1$  as the low-density phase,  $\alpha/\beta = 1$  as the coexistence phase, and  $\alpha/\beta > 1$ <sup>231</sup> as the high-density phase. This is consistent with the known results obtained using exact <sup>232</sup> methods in [16]. We depict our perturbative phase diagram of the open ASEP in Figure 2.

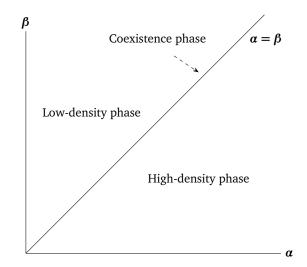


Figure 2: The phase diagram of the open ASEP. The horizontal axis represents the adsorption rate at site i = 1, while the vertical, the desorption rate at site i = L. This recovers the perturbative part of the known phase diagram of the open ASEP, obtained exactly in [16].

#### 233 4.2 Phase diagram of the open ASEP-LK

We can also compute the stationary state of the open ASEP-LK by turning on a and b. Just as in the case of the open ASEP, we have

$$A_{i} = \frac{a(i-1) + aq^{i} + (\alpha - a)q^{L-i} + (\alpha - \alpha - ai)q^{L}}{1 - a^{L}},$$
(4.22)

$$B_{i} = \frac{b(i-1) + \beta + (b-\beta)q^{i} - bq^{L-i} + (b-bi)q^{L}}{1 - q^{L}},$$
(4.23)

from which we can compute the stationary state of the open ASEP-LK at leading order in  $\epsilon$ . In particular at q = 0,  $p_N$  can be expressed concisely as

$$p_N = \left(-\frac{a}{b}\right)^n \frac{\left(-L - \frac{a}{a} + 1\right)_n}{\left(\frac{\beta}{b}\right)_n},\tag{4.24}$$

where  $(x)_n \equiv \prod_{i=0}^{n-1} (x+i)$  is the Pochhammer symbol. One can then compute  $\langle \tau \rangle_i = \sum_{k=1}^{L} \frac{p_N}{\sum_{k=1}^{L} p_N} \sum_{k=0}^{L} \frac{p_N}{p_N}$  and express it using hypergeometric functions, but we will not discuss this further as it will just be unnecessarily complex. We plot  $\langle \tau \rangle_i$  for some parameters in Figure 3.

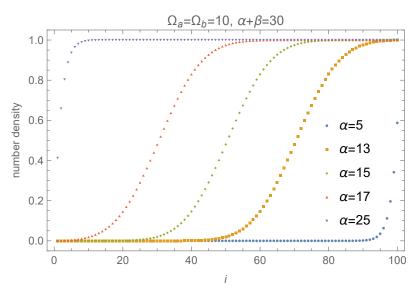


Figure 3: Plot of the particle number densities of open ASEP-LK (at q = 0) as functions of lattice sites *i*. We take the number of lattice sites to be L = 100. We set  $\Omega_a \equiv aL = 10$ ,  $\Omega_b \equiv bL = 10$  and varied  $\alpha$ ,  $\beta$  while fixing  $\alpha + \beta = 30$ . We sampled  $\alpha = 5$ , 13, 15, 17, 25 in the plot. We see that  $\alpha = 5$  is in the low-density phase,  $\alpha = 13$ , 15, 17, the domain-wall phase, and  $\alpha = 25$ , the high-density phase, consistent with analytic computations.

We now take the thermodynamic limit,  $L \to \infty$ . For the sake of manageability we will only consider the bulk region of the open chain, so that we take  $i \to \infty$  at the same time while fixing  $x \equiv i/L$ . We will also take  $a, b \to 0$  while fixing  $\Omega_a \equiv aL$  and  $\Omega_b \equiv bL$  – otherwise the collective effect of the bulk ad/desorption will dominate the physics and there will be no interesting phase transitions.

Let us compute  $\langle \tau \rangle_i = \sum_{N=L+1-i}^{L} p_N / \sum_{N=0}^{L} p_N$ . At large *L* and at fixed *x*,  $\Omega_a$ ,  $\Omega_b$ , it simply becomes

$$\rho(x) \equiv \langle \tau \rangle_i = \begin{cases} 1 & \text{when } p_{L-i+1}/p_{L-i} > 1 \\ 0 & \text{when } p_{L-i+1}/p_{L-i} < 1 \end{cases},$$
(4.25)

248 where we have

$$\frac{p_{L-i+1}}{p_{L-i}} = \frac{\Omega_a x + \alpha}{\Omega_b (1-x) + \beta} + O(L^{-1}), \tag{4.26}$$

for general 0 < q < 1. This means that the domain-wall that separates the low- and the highdensity phase happens at  $x_d$  (the former appears for  $x < x_d$  and the latter,  $x > x_d$ ), given 251 by

$$x_d = \frac{\Omega_b - \alpha + \beta}{\Omega_a + \Omega_b}.$$
(4.27)

We call this the domain-wall phase (called the shock phase in [12]).<sup>3</sup> Additionally, when  $x_d > 1$ , the system is in the low-density phase, whereas when  $x_d < 0$ , it is in the high-density phase. Summarising this, we have the low-density phase when  $\beta > \alpha + \Omega_a$ , the domainwall phase when  $\alpha - \Omega_b < \beta < \alpha + \Omega_a$ , and the high-density phase when  $\beta < \alpha - \Omega_b$ . This is consistent with the results obtained using the (theoretically unjustified but numerically confirmed) mean-field approximation in [12]. We depict our perturbative phase diagram of the open ASEP-LK in Figure 4.

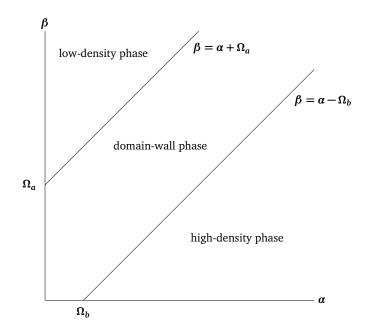


Figure 4: The phase diagram of the open ASEP-LK. The horizontal axis represents the adsorption rate at site i = 1, while the vertical, the desorption rate at site i = L. This recovers the perturbative part of the known phase diagram of the open ASEP-LK, obtained using the mean-field approximation in [12].

## **5** Discussions and outlook

In this paper, we studied the effect of perturbation on generic closed exclusion processes. We first derived the formula that expresses the stationary state of closed processes (infinitesimally) perturbed by ad/desorption in terms of that of the unperturbed system. The rates of ad/desorption did not have to be homogeneous in sites, so as a consequence we proved the formula in [11] while generalising it. We pointed out that our formula is a result of the simple degenerate perturbation theory on non-normal matrices.

As an application of the formula, we drew the perturbative part of the phase diagram of the open ASEP(-LK), which agreed with known results. For the open ASEP we recognised three distinct phases, called the low-density, the coexistence, and the high-density phases. For the

<sup>&</sup>lt;sup>3</sup>The position of the domain wall  $x_d$  is indeed consistent with numerics, see Figure 3. We expect the position to lie at i = 70, 50, 30 for a = 13, 15, 17, respectively.

open ASEP-LK, on the other hand, we recognised the low-density and the high-density phases,
as well as the domain-wall phase in which the system contains a domain wall separating the
low- and the high-density regions. It is important that these results were obtained without
using any theoretically unjustified approximations – we exactly know when and how much
our approximation breaks down.

There are a number of interesting future directions. First of all, it would be interesting to continue the perturbation theory to higher orders in  $\epsilon$ . For example, if we compare the phase diagram of [12] with ours, we notice that the phase boundaries are not exactly straight, *i.e.*,  $\beta$  at the critical value is not a linear function of  $\alpha$ . It would be beneficial to compute the form of the phase boundaries at higher orders in perturbation theory to explain this.

Secondly, it would be interesting to apply our method to other systems of interest. For 279 example, it would be interesting to apply it to the multi-lane ASEP [17] or to the ASEP(-280 LK) with inhomogeneous hopping rates [18]<sup>4</sup>. It would also be interesting to study the open 281 ASEP-LK by starting the perturbation from the exactly known stationary state of the open ASEP. 282 Note that what we would need to do is in general non-degenerate perturbation theory. The 283 result would allow us draw wider region of the phase diagram upon taking the thermodynamic 284 limit. In particular, observing the three-phase coexistence predicted in [12] would be very 285 interesting. 286

Studying the relaxation dynamics in perturbation theory is also interesting. One could, for example, compute the low-lying spectra and the corresponding states for the same class of theories at leading order in perturbation theory. In fact, [11] conjectures such a formula for the closed ASEP-LK, so it would be interesting to start by proving it.

It would be important to justify the mean-field approximation theoretically as well. One 291 could for example compute the two-point functions perturbatively in  $\epsilon$ ; If they factorise in 292 the thermodynamic limit, the mean-field approximation is justified at least perturbatively. It 293 would also be useful to justify it without relying on other perturbation theory at all. In this con-294 text, it might be worthwhile to rewrite the open ASEP-LK in the language of one-dimensional 295 (non-Hermitian) spin chains. The mean-field approximation can then be justified when the 296 model flows to the free fixed-point in the infrared. It would also be interesting to come up 297 with a model which is strongly-coupled in the infrared, where the mean-field approximation 298 cannot be justified. Incidentally, in terms of the field theoretic understanding of the exclusion 299 processes, interpreting the asymmetric hopping parameter q as an imaginary vector potential 300 is also interesting [21]. Because the  $q \rightarrow 0$  limit corresponds to the limit of large imaginary 301 vector potential, one might be able to use effective field theory to study such regions [22-28]. 302 Lastly, studying the relationship between the general solvable exclusion processes with 303 other models with  $U_q(sl(2))$  symmetry would be interesting. In particular, the SYK model (a 304 quantum mechanical model with all-to-all random interactions of N fermions) in the double-305 scaling limit [29-31] is known to possess such a symmetry and it would be interesting to 306 connect them further. It would also be interesting to interpret it in terms of Jackiw-Teitelboim 307 gravity [32, 33], which is believe to be dual to the SYK model in the context of the AdS/CFT 308 correspondence [34]. 309

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<sup>&</sup>lt;sup>4</sup>Studying the latter in relation to sine-square deformation and other similar deformations [19, 20] might be also interesting.

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