



A FINITE ELEMENT CODE FOR GEOTECHNICAL SIMULATIONS

INPUTS FROM PUBLICATIONS

-

SERIES A: VIBRATORY PILE DRIVING IN WATER-SATURATED
SAND

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The user routines referred to in the following are for the Ubuntu version of `numgeo`. The data types are slightly different for the Windows version. The user routines for the Windows are contained in the folder 'input_windows' shipped with this document.

1 Vibratory pile driving in water-saturated sand

Figure 1: Spatial distribution of pore water pressure during vibratory driving

In this series the simulation of vibratory pile driving in water-saturated sand according to the simulations reported in [1] is explained. The input file shipped together with this document and further described in the following is identical to the one used in [1] from a technical point of view. Note that the input files used in [2] are very similar to the one explained in this document.

All information about the numerical model used for the analysis can be obtained from [1]. This document only explains the definitions in the input file and not the general set-up of the numerical model.



The simulation of the input file adopted here has been performed with the corresponding release of `numgeo` (August 2023). Using older releases, you might have to adjust the input file in order to be able to perform the simulation since keyword definitions have changed. In any case, using the latest release is recommended.

1.1 Input file

The input file is split into two files: the mesh (`mesh.inp`) and the step file (`step.inp`). First, the mesh file is briefly explained.

1.1.1 Mesh file

`u8-solid-ax` elements are used for the pile and the unbalances, which are placed on top of the pile and are connected to it by spring elements. For the soil, axisymmetric `u-p` finite elements with 8 nodes discretising the displacement \mathbf{u} of the solid phase and 4 nodes discretising the pore water pressure p^w are utilised (termed `u8p4-sat-ax-red` element). A reduced integration scheme with 4 integration points is adopted. The corresponding definition can be found in line 7550 of `mesh.inp`.

```

0 .
1 .
2 *Element , type=u8p4-sat-ax-red
3 .
4 .

```

Hence, quadratic interpolation functions are used for the solid displacement while linear interpolation functions are applied for the pore water pressure (i.e. Taylor-Hood element formulation [4]).

1.1.2 Properties, initial conditions and step definitions (Step file)

The step file (*step.inp*) first includes the mesh file by defining:

```
0 *include = mesh
```

In the model test by J. Vogelsang [6], two pore water pressure transducers recorded the development of pore water pressure at different spatial positions during driving. In order to compare the measurements with the results of the simulations, the pore water pressure is also recorded in the numerical analysis. The points are defined by specifying their spatial position:

```
0 *nset, find, nset = top-pw
1 0.0386,0.6,0.0
2
3 *nset, find, nset = bottom_pw
4 0.025,0.35,0.0
```

The penalty method is used to enforce the contact constraints as is given in Listing 1. No separation between soil and pile is possible once the contact is active. The penalty factor is $3 \cdot 10^6$ kN/m³. A simple Coulomb friction model is adopted, which properties are defined in a user-file, which is explained later in Section 1.2.1. The contact between soil and pile is discretised using the element-based mortar contact discretisation technique (line 5), which implementation is discussed in detail in [3]. A small value for the contact clearance is defined, which leads to a small value of Initial contact pressure. This is beneficial since the Initial stress state is closer to equilibrium, in which the earth pressure is acting as contact pressure.

```
0 **----- contact -----
1
2 *interaction, name = penalty, mechanical = penalty, no separation, user
3 3d6
4 *friction, model = mc
5 *contact pair, interaction = penalty, discretisation = element mortar
6 soil_pile_all, pile_soil_all
7 *contact options, name = penalty, clearance = 2d-5
```

Listing 1: Definition of the contact

Then the definitions of the solid sections and material definitions are given. The soil is modelled with the material *hypo_soil* while pile and unbalance are modelled elastically. However, the elastic properties are set arbitrarily high since no self-deformations occurred in the tests. Pile and unbalance also don't have a density, since their self-weight and inertia will be applied by concentrated forces and mass points, respectively.

```
0 **----- solid section -----
1
2 *solid section, elset = soil_all, material = hypo_soil
3 *solid section, elset = pile_all, material = elastic
4 *solid section, elset = unbalance, material = elastic
5
6 *material, name = elastic, phases = 1
7 *mechanical = linear_elasticity
8 5d10, 0.0
9 *density
10 1d-10
```

Only the hypoplastic model with intergranular strain extension is considered here as constitutive model for the soil. The input file for the simulations using Sanisand is easily obtained by replacing the following definitions. The parameters of the hypoplastic model with intergranular strain extension according to [1] are given in lines 2-3 of Listing 2.

An artificial viscosity is used to achieve better convergence. The adopted values are further described in [1]. To avoid positive mean effective stresses (in mechanical sign convention), a minimum value is set in line 10-11.

Two densities have to be given which are the density of the solid grains and the density of water. In addition, the bulk modulus of the pore water and the permeability have to be defined. In *numgeo*, the permeability equals the dynamic viscosity (set to 10^{-6} kPas) multiplied with the hydraulic conductivity k^w and divided by the unit weight of water ($\gamma_w = 10$ kN/m³). As explained in [1] (and further evaluated in [2]) the Kozeny-Carman relation is used to consider the influence of the porosity on the permeability.

```
0 *material, name = hypo_soil, phases = 2
1 *mechanical = hypoplasticity
2 0.577, 0.0, 19d6, 0.285, 0.549, 0.851, 1.15, 0.1
```

```

3 0.32, 1.2, 2.4, 5d-5, 0.08, 7, 0.
4
5 *mechanical viscosity = linear
6 0.5, 2.0, 0.1, 1, 0.1, 1
7 *minpressure
8 -0.01
9
10 *density
11 2.65, 1.0
12
13 *bulk modulus
14 1.2d4
15 *permeability = kozeny-carman
16 308, 0.5d-3
17 *dynamicviscosity
18 1.0d-6

```

Listing 2: Definition of the soil material

The initial conditions for the state variables are defined in user subroutines (see the files shipped with this document and the explanations given later in Section 1.2.1). The void ratio used for the calculation of the density of the soil is given in line 3. This value does not necessary have to coincide with the value given in the user initial state file. The initial stresses according to the values given in [1] are defined in lines 8-11. A hydrostatic pore water pressure is defined in lines 13-14 for the entire soil mass.

```

0 **----- Initial conditions -----
1
2 *initial conditions, type = void ratio, default
3 soil_all, 0.637
4
5 *initial conditions, type = state variables, user
6 soil_all
7
8 *initial conditions, type = stress, geostatic
9 soil_all, 0.81, -0.1, 0, -8.266, 0.455, 0.455
10 pile_all, 1, 0, 0, 0, 0.5, 0.50
11 unbalance, 1, 0, 0, 0, 0.5, 0.50
12
13 *initial conditions, type = pore water pressure, default
14 soil_all, 0.0, 8.1, 0.81, 0.0

```

For the analysis, different time distributions of boundary and loading conditions have to be defined. This is done using the following amplitude definitions:

```

0 **----- define amplitudes -----
1 *amplitude, name = loadingramp, type = ramp
2 0.0, 0.3, 1.0, 1.00
3
4 *amplitude, name = risingsine, type = rising_sine
5 0.075, 157.0, 1.0

```

The amplitude *loadingramp* defines a linear increase of a quantity beginning with the relative value 0.3 at the time $t_1 = 0$ and the relative value 1 at the time $t_2 = 1$. To apply the cyclic loading by the unbalances, a sinusoidal amplitude with a frequency of 25 Hz is defined in addition. For this amplitude, a linear increase for the first 0.075 s is assumed, considering that in the model tests, the unbalances also required a short time to achieve the targeted frequency.

1.1.3 Steps and loading definitions

The first step is defined in Listing 3. Line 3 starts the step environment and defines the name of the step. The analysis type of the step is given in line 4, which will be a geostatic step considering non-linear geometry effects (*nlgeom*). The keyword *Body force* imposes a gravitational force which is applied instantaneous (*Instant*). The element set *soil_all* is loaded by the gravity (Amplitude 10 m/s², directed downwards with the normalized vector of the gravity $\vec{\mathbf{b}} = \{0, -1, 0\}$).

The boundary conditions are specified from line 9 to line 14. All nodes of the pile, unbalance and the soil are constraint in 1- and 2-direction (i.e. radial and vertical direction). Therefore, no displacement is possible for any node in the model. The pore water pressure is prescribed hydrostatically using a boundary condition of type *hydrostatic*. A small stabilising pressure of 0.1 kPa is applied at the top surface of the soil.

Not active in the step but already being defined for the next step is a connector element, which connects unbalances and pile. The connector element mimics the spring between unbalances and pile used to measure the force. The spring stiffness is 173000 kN/m.

The output demand is specified from line 32 to line 44. The output is written in the vtk format (suitable for ParaView) and is of ASCII type. The node output includes the displacement (**u**) and pore water pressure (**pw**) of the nodes. The element output includes the stress (**s**), strain (**e**) as well as the void ratio (**void_ratio**) and the contact output variables (**contact**). The contact output contains contact stresses as well as contact distances for each contact node. In addition, a print output is defined in lines 32-44. The displacement and the pore water pressure is printed for several nodes of the model. In addition, the connector force of the spring between unbalances and pile is written and the acceleration. From lines 41-42 the displacement and coordinates of all nodes of the soil are defined but not active. This output definition creates considerable data as is required to create the incremental displacement fields presented in [1].

```

0  **----- steps -----
1
2  *step, name = step1, inc = 1, nlgeom
3  *geostatic
4
5  *body force, instant
6  soil_all, grav, 10, 0, -1, 0
7
8  *boundary
9  pile_all, u1, 0.0
10 soil_all, u2, 0.0
11 soil_all, u1, 0.0
12 pile_all, u2, 0.0
13 unbalance, u1, 0.0
14 unbalance, u2, 0.0
15
16 *boundary, type = hydrostatic
17 soil_all, pw, 10.0, 0.81
18 *dload, instant
19 _soil_top_s3, p3, -0.1
20 *dload, instant
21 _soil_top_s4, p4, -0.1
22
23 *connector element, instant
24 unbalance_con, pfahl_con, u2, 173000
25
26 *output, field, vtk, ascii
27 *node output, nset = soil_all
28 u, pw
29 *element output, elset = soil_all
30 s, e, contact, void
31
32 *output, print
33 *node output, nset = u_oscillator
34 u
35 *node output, nset = unbalance_con
36 con_f
37 *node output, nset = top_pw
38 pw
39 *node output, nset = bottom_pw
40 pw
41 **node output, nset = soil_all
42 **u, coords
43 *node output, nset = pile_acceleration
44 a
45
46 *end step

```

Listing 3: Definition of the first step

In the second step given in Listing 4, the correct kinematic boundary conditions for the soil are established. The bottom of the soil is constraint in the vertical direction in line 19. The back-side is constraint in the radial direction in line 20. The unbalance is fixed in space in this step. The output remains identical to the first step.

```

0  **----- steps -----
1
2  *step, name = step2, inc = 100000, nlgeom

```

```

3 *static
4 1,1.0,0.001,1
5
6 *body force , instant
7 soil_all , grav , 10.0, 0, -1, 0
8
9 *dload , instant
10 _soil_top_s3 , p3 , -0.1
11 *dload , instant
12 _soil_top_s4 , p4 , -0.1
13
14 *connector element
15 unbalance_con , pfahl_con , u2 , 173000
16
17 *boundary
18 pile_all , u1 , 0.0
19 soil_bottom , u2 , 0.0
20 soil_back , u1 , 0.0
21 unbalance , u1 , 0.0
22 unbalance , u2 , 0.0
23
24 *boundary , type = hydrostatic
25 soil_all , pw , 10.0 , 0.81
26
27 *output , field , vtk , ascii
28 *node output , nset = soil_all
29 u , pw
30 *element output , elset = soil_all
31 s , e , contact , void
32
33 *output , print
34 *node output , nset = u_oscillator
35 u
36 *node output , nset = unbalance_con
37 con_f
38 *node output , nset = top_pw
39 pw
40 *node output , nset = bottom_pw
41 pw
42 **node output , nset = soil_all
43 **u , coords
44 *node output , nset = pile_acceleration
45 a
46
47 *end step

```

Listing 4: Definition of the second step

In the third step given in Listing 5 the unbalance is release and the weight of unbalance and pile applied (lines 26-29).

```

0 **----- steps -----
1
2 *step , name = step3 , inc = 100000 , nlgeom
3 *static
4 0.001,1.0,0.000001,0.02
5
6 *body force , instant
7 soil_all , grav , 10, 0, -1, 0
8
9 *dload , instant
10 _soil_top_s3 , p3 , -0.1
11 *dload , instant
12 _soil_top_s4 , p4 , -0.1
13
14 *connector element
15 unbalance_con , pfahl_con , u2 , 173000
16
17 *boundary
18 pile_all , u1 , 0.0
19 soil_bottom , u2 , 0.0
20 soil_back , u1 , 0.0
21 unbalance , u1 , 0.0

```

```

22
23 *boundary, type = hydrostatic
24 soil_all, pw, 10.0, 0.81
25
26 *cload, amplitude = loadingramp
27 unbalance_con, 2, -0.1119
28 *cload, amplitude = loadingramp
29 pfahl_con, 2, -0.0218
30
31 *controls, global, deactivate
32 *controls, u, modify
33 0.05, 0.05, 0.05, 4e-9, 4e-13, 2
34
35 *output, field, vtk, ascii
36 *frequency=20
37 *node output, nset = soil_all
38 u, pw, con_f
39 *element output, elset = soil_all
40 s, e, contact, void
41
42
43 *output, print
44 *node output, nset = u_oscillator
45 u
46 *node output, nset = unbalance_con
47 con_f
48 *node output, nset = top_pw
49 pw
50 *node output, nset = bottom_pw
51 pw
52 **node output, nset = soil_all
53 **u, coords
54 *node output, nset = pile_acceleration
55 a
56
57 *end step

```

Listing 5: Definition of the third step

The fourth step the simulation of the vibratory driving process is simulated. A dynamic analysis is performed. 6.35 s of driving are simulated according to line 4. A dynamic analysis requires a time integration scheme (since the displacement is discretised, the velocity and the accelerations are unknown and have to be determined). The Hilber-Hughes-Taylor (HHT) time integration scheme is chosen here and a slight numerical damping with $\alpha = -0.05$ is defined. Since inertia effects have to be accounted for, the self-weight of pile and unbalances is considered by defining two mass points (lines (18-21)). The loading by the unbalances is defined in lines 34-35 using the previously defined amplitude.

```

0  **----- steps -----
1
2 *step, name = step4, inc = 1000000, nlgeom
3 *dynamic
4 0.001, 6.35, 0.000001, 0.001
5 *hilber-hughes-taylor=-0.05
6
7 *body force, instant
8 soil_all, grav, 10, 0, -1, 0
9
10 *dload, instant
11 _soil_top_s3, p3, -0.1
12 *dload, instant
13 _soil_top_s4, p4, -0.1
14
15 *connector element
16 unbalance_con, pfahl_con, u2, 173000
17
18 *mass point, instant
19 unbalance_con, 2, 0.01119
20 *mass point, instant
21 pfahl_con, 2, 0.00218
22
23 *boundary
24 pile_all, u1, 0.0

```

```

25 soil_bottom , u2, 0.0
26 soil_back , u1, 0.0
27 unbalance , u1, 0.0
28 soil_top , pw, 0.0
29
30 *cload , instant
31 unbalance_con , 2, -0.1119
32 *cload , instant
33 pfahl_con , 2, -0.0218
34 *cload , amplitude = risingsine
35 load , 2, -0.2329845
36
37 *controls , global , deactivate
38 *controls , u, modify
39 0.05, 0.05, 0.05, 4e-9, 4e-13, 2
40 *controls , pw, deactivate
41
42 *output , field , vtk , ascii
43 *frequency = 10
44 *node output , nset = soil_all
45 u, v, a, pw, con_f
46 *element output , elset = soil_all
47 s, contact , void , darcs_w1 , darcy_w2
48
49 *output , print
50 *node output , nset = u_oscillator
51 u
52 *node output , nset = unbalance_con
53 con_f
54 *node output , nset = top_pw
55 pw
56 *node output , nset = bottom_pw
57 pw
58 **node output , nset = soil_all
59 **u, coords
60 *node output , nset = pile_acceleration
61 a
62
63 *end step

```

Listing 6: Definition of the fourth step

In the last step, which is skipped here, the driving force is deactivated and the fading out of the pile movement is studied.

1.2 User files

The user files used to define the contact properties and the initial state variables of the soil are explained in more detail in the following.

1.2.1 User contact properties

As explained earlier, the contact properties are defined by a user subroutine. This is necessary, since the contact surface is defined for the entire pile length (including the thin extension for the zipper technique, see [1]). However, friction is only to be considered at the tip and the shaft up to the ground surface. Therefore, using the user subroutine `user_contact_properties`, the friction coefficient is defined coordinate-dependent. Of course, it would alternatively also be possible to define multiple surface pairs with different friction coefficients instead.

As can be seen from the `user_contact_properties` fortran file, friction is only considered at the tip and active shaft area of the pile, which also changes with displacement of the pile in order to consider the ongoing penetration.

```

0 subroutine user_contact_properties(istep , node , slave , nprops , interaction_type , step_time , coords , &
1 coords_connected , disp , disp_connected , props) bind(c, name='user_contact_properties')
2 use, intrinsic :: iso_c_binding
3 implicit none
4 integer(c_int) , intent(in) :: istep
5 integer(c_int) , intent(in) :: node
6 logical , intent(in) :: slave
7 integer(c_int) , intent(in) :: nprops

```

```

8 character(c_char) , intent(in) :: interaction_type(*)
9 real(c_double) , intent(in) :: step_time
10 real(c_double), dimension(3) , intent(in) :: coords
11 real(c_double), dimension(3) , intent(in) :: coords_connected
12 real(c_double), dimension(3) , intent(in) :: disp
13 real(c_double), dimension(3) , intent(in) :: disp_connected
14 real(c_double), dimension(nprops) , intent(inout) :: props
15
16 props(:) = 0.0d0
17 props(1) = 3d3 ! tangential stiffness
18 if(slave .eqv. .false.) then
19   if(coords(2) + disp(2) > 0.641 + disp(2) .and. coords(2) < 0.8138 - disp(2) ) then
20     props(2) = 0.194d0 ! friction coefficient
21   endif
22 else
23   if(coords_connected(2)+ disp_connected(2) > 0.641 + disp_connected(2)) then
24     props(2) = 0.194d0 ! friction coefficient
25   endif
26 endif
27
28 end subroutine user_contact_properties

```

1.2.2 User initial state variables properties

Because relative density is stress dependent in the hypoplastic model, the initial void ratio is initialized considering the decrease with increasing mean effective stress using Bauer's formula. The void ratio (statev(1)) and the intergranular strain in vertical direction (statev(3)) are initialized using the the `user_initial_state_variables` fortran file. Detailed explanations may be found in [1].

```

0 subroutine user_initial_state_variables(ie, igp, ndim, nstatev, material, coords, statev) bind(c, name
   = 'user_initial_state_variables')
1 use, intrinsic :: iso_c_binding
2 implicit none
3 integer(c_int) , intent(in) :: ie
4 integer(c_int) , intent(in) :: igp
5 integer(c_int) , intent(in) :: ndim
6 integer(c_int) , intent(in) :: nstatev
7 character(c_char) , intent(in) :: material(*)
8 real(c_double), dimension(3) , intent(in) :: coords
9 real(c_double), dimension(nstatev), intent(inout) :: statev
10
11 real(c_double) :: h_top, void0, K0, surface_load, n, gamma_tot, pw, stress, trT, hs, n_hyp
12
13 statev(:) = 0.0d0
14
15 h_top = 0.81d0
16 void0 = 0.637d0
17 K0 = 1.0d0 - sin(0.577d0)
18 hs = 19000000.0d0
19 n_hyp = 0.285
20
21 ! external load on top surface
22 surface_load = 0.1d0
23
24 ! total unit weight
25 n = void0 / (1.0d0 + void0)
26 gamma_tot = ((1.0d0 - n) * 2.65d0 + n) * 10.0d0
27
28 ! pore water pressure
29 pw = (h_top - coords(2)) * 10.0d0
30
31 ! stress (vertical and mean effective stress)
32 stress = gamma_tot * (h_top - coords(2)) + surface_load - pw
33 trT = (1.0d0 + 2.0d0 * K0) * stress / 3.0d0
34
35 statev(1) = void0 * exp(-(trT/hs) ** n_hyp)
36
37 statev(1+2) = -0.00005
38
39 end subroutine user_initial_state_variables

```

1.3 Results of the simulation

During the calculation or after the it is finished (the command window is ready for a new command), open the .sta file first and check that the simulation was successful by identifying that all steps have been completed successfully. If the calculation immediately stops, check the error message in the .log file and if any error files were generated in the calculation folder.

The pile penetration vs. time of vibration is depicted in Fig. 2 for the simulations and the measurements by J. Vogelsang [5, 6].

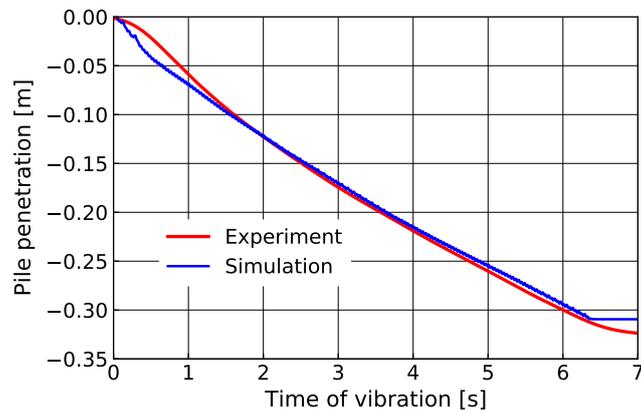


Figure 2: Pile penetration vs. time of vibration

References

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