

HYPACT

HYbrid Particle And Concentration Transport Model

Version 1.5

User's Guide

by

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About *HYPACT*

The *HYPACT* (HYbrid Particle And Concentration Transport) model code has been developed to simulate the motion of atmospheric tracers under the influence of atmospheric flow, including turbulence. In its current form, it is set up to be driven by meteorological output from *RAMS* (Regional Atmospheric Modeling System), but the basic design allows the flexibility, with minor code changes, to provide a variety of other inputs as well, including analyzed, observational data.

Although *RAMS* can itself perform a similar function by simulating the motions of any number of Eulerian tracer fields, *HYPACT* has certain advantages because of the ability to run the dispersion as a post-processing step and because of its Lagrangian formulation. The advantage is greatest near a source region for the tracer when the source is of small scale and unresolved on the Eulerian grid. *RAMS* would necessarily represent the source by a volume no smaller than one grid cell, and would immediately begin diffusing the tracer to the adjacent cells. A Lagrangian model, on the other hand, is capable of representing a source of any size, and of maintaining a concentrated, narrow tracer plume downwind of the source until atmospheric dispersion dictates that the plume should broaden. An important feature of *HYPACT* is the “hybrid” capabilities: once a plume of Lagrangian tracer particles becomes sufficiently broad downstream from a source region, it can be converted to a resolved concentration field and advected using an Eulerian formulation. This hybrid approach allows high particle emission rates to be specified for a source to achieve good plume resolution, without retaining all particles so far into the future that excessively large numbers of them accumulate. Tracer sources that are well resolved on a grid may alternatively be specified as gridded sources in *HYPACT* and transported entirely by the Eulerian formulation. *HYPACT* carries out Eulerian tracer prediction much faster than *RAMS* because it does not predict its own wind, pressure, temperature, moisture, or turbulence fields.

In order to run *HYPACT*, *RAMS* must be run first to generate a series of output “analysis” files. These contain wind, potential temperature, and turbulent kinetic energy fields at the output times of the atmospheric simulation. *HYPACT* reads these data from the analysis files and interpolates them in time between file times (for each *HYPACT* timestep) for determining transport and diffusion. For Eulerian concentration prediction, *HYPACT* assigns the time-interpolated data to a series of nested grids identical with those of the *RAMS* simulation that generated the analysis files. Concentrations are predicted on these grids given user-specified initial values and source characteristics. For Lagrangian particle prediction, the time-interpolated wind and turbulence data are also interpolated in space to the location of each Lagrangian tracer particle. If nested grids are used in the *RAMS* simulation, the finest grid data defined at the location of each particle is used for the interpolation. The particles are moved through space and time based on the interpolated wind velocity plus a random motion scaled to the local turbulent intensity. A gravitational settling velocity may also be superimposed on the particle motion.

HYPACT sources, species and emission can either be defined in the *HYPACT* namelists, or in database files. This is designed to allow both quick test runs, and a range of sophisticated scenarios from common data.

In addition to Eulerian concentration field outputs on the *RAMS* grids, concentration can be periodically diagnosed from the particle distribution on the *RAMS* grids (assuming a specified mass represented by each particle), and the combined Eulerian and Lagrangian concentration fields computed and output. The concentration is useful for contouring plume morphologies, determining instantaneous or time-integrated dosages at specified locations, computing chemical reaction rates between different tracer fields, etc. The raw particle files can also be output, allowing the user to diagnose the particle concentrations on a finer concentration grid.

HYPACT version 1.5 is FORTRAN 90 compliant. UNIX/Linux make commands and a system of make files and makefile includes are used for compiling the code. Similar to *RAMS*, *HYPACT* is in a state of

continual evolution. As with the *RAMS* documentation, we have split the User's Guide into multiple sections. This document will describe the namelist configuration parameters. Other documents will cover the building, installation, and execution.

New in HYPACT Version 1.5

New features since *HYPACT* version 1.2 include:

- Further modifications to the code structure and memory allocation, following the use of Fortran 90 user-defined data types as was done in *RAMS* v5. All include files have been converted to F90 modules.
- Supports *RAMS* v5 and *REMU* v2.5
- Backwards trajectories
- New way to specify namelist times
- New feature for *RAMS* v5 compatibility to allow nested grid analysis files to be written at different frequencies

Note that v1.3 and 1.4 were temporary developmental versions that were not officially released.

New in HYPACT Version 1.2

New features since *HYPACT* version 1.1 include:

- Modified code structure and memory allocation. As a result, this version of *HYPACT* is far more robust than previous version. We strongly recommend users of previous versions upgrading to this version, as we will no longer support the older versions. Specifically, many of the routines now use "implicit none" statements and all `c iralloc` memory allocations are now done with FORTRAN 90 allocations. This has allowed us to do bounds checking on arrays which in turn has resulted in the cleaning up of a number of known and unknown bugs.
- Several long time bug fixes and *RAMS* / *HYPACT* code consistency checks.
- A fix to the hybrid conversion routine, ensuring stable plumes of Lagrangian particles advecting (about) 45 degrees from an north, south, east or west directions do not prematurely get converted to Eulerian concentrations.
- Accepts 1 argument, `-f`, for pointing to non-standard [*HYPACT_IN*](#) file names, i.e.

```
hypact-1.2.0 -f <namelist file>
```

HYPACT Configuration Parameters

Following are some *HYPACT* parameters that are set at compilation time in *hyp_coms.f90*. They are used to dimension several arrays, primarily in the *hyp_coms* module. These are maximum settings of these variables and are defaulted to rather large values. These settings do not affect memory usage much, as all significant memory is dynamically allocated. The only reason to change these values is if you have run an extremely large RAMS simulation and/or have many nested grids.

maxg integer	maxg must be greater than or equal to the number of grids used in the <i>RAMS</i> simulation used to drive <i>HYPACT</i> .
maxi maxj maxk integer	must be greater than or equal to the largest number of grid points spanning the x-direction (east to west), y-direction (south to north) and z-direction (vertical), respectively, of any grid used in a <i>RAMS</i> simulation used to drive <i>HYPACT</i> .
maxhi maxhj maxhk integer	These variables specify upper bounds on the number of grid points spanning the i, j, and k directions, respectively, of the special grid defined in <i>HYPACT</i> (<i>h-grid</i>). The <i>h-grid</i> is a composite of all nested grids used in a <i>RAMS</i> simulation used to drive <i>HYPACT</i> . It has variable spatial resolution according to the mesh sizes, overall dimensions, and locations of all grids in <i>RAMS</i> . The <i>h-grid's</i> mesh size is always at least as fine as, and often finer than, the finest <i>RAMS</i> grid mesh size at any given location. The required number of points spanning each direction of the <i>h-grid</i> is computed by a complicated algorithm, but is usually less than the sum over all <i>RAMS</i> grids of the numbers of points spanning each corresponding direction of the <i>RAMS</i> grids. Thus, if for example <i>RAMS</i> uses 3 grids each of (i,j,k) dimensions (30,40,50), the <i>h-grid</i> would likely have dimensions somewhat smaller than (90,120,150). These parameters need not match the exact size of the <i>h-grid</i> ; they only have to be large enough to contain it.
maxfiles integer	maxfiles must be greater than or equal to the largest number of <i>RAMS</i> “analysis” files used in a <i>HYPACT</i> simulation that have the path and filename prefix given by hyppref (not just those within the defined <i>HYPACT</i> simulation time).
maxspec integer	maxspec must be greater than or equal to the largest number of species to be used in a <i>HYPACT</i> simulation.
maxsrc integer	maxsrc must be greater than or equal to the largest number of sources to be used in a <i>HYPACT</i> simulation.
maxim integer	maxem must be greater than or equal to the largest number of emissions to be used in a <i>HYPACT</i> simulation. The number of emissions used is the tally of emissions turned on by \$EMISSIONS namelist variable iemit .

--	--

polypts integer	polypts must be greater than or equal to the largest number of vertices used to define any polygon emission source in a <i>HYPACT</i> simulation. The number of vertices used for each polygon is specified by the \$SOURCES namelist variables polylat and polylon .
maxent integer	maxent specifies the size of a table of normally-distributed numbers used in <i>HYPACT</i> . A value of approximately 1000 provides adequate density in the table.

HYPACT Namelist and Database Files

The user has the flexibility of emitting particles and concentrations from different locations, at different times, at different rates, and with different source region characteristics. Each particle carries with it a label indicating the source from which it was emitted, so the particles can be identified accordingly throughout their journey.

A database approach is used in the specification of species and sources in the *HYPACT_IN* namelist file. As such, any number of species and sources may be individually defined in the [\\$SPECIES](#) and [\\$SOURCES](#) namelists (within the limits of the *HYPACT* parameter settings). Any combinations of species and sources can then be specified in conjunction with emission data (rate, duration, etc) in the [\\$EMISSIONS](#) namelist. Thus, the *HYPACT_IN* namelist file can contain a comprehensive species and sources datasets, with the [\\$EMISSIONS](#) namelist using only that data required for the current run. In addition to this flexibility, any number of emissions (within the limits of the *HYPACT* parameter settings) can be specified in the [\\$EMISSIONS](#) namelist.

Alternatively, the species, source and emissions specifications included in the namelists can be carried in separate “database” files. If using species, source and emissions database files, only minimal number of parameters are required in the [\\$SPECIES](#), [\\$SOURCES](#) and [\\$EMISSIONS](#) namelists. The format for these files is discussed in the following sections.

Two other namelists [\\$GENERAL](#) and [\\$OUTPUT](#) control the other aspects of running the model.

\$GENERAL Namelist

<p>ihyprun</p> <p>integer</p>	<p>The <i>HYPACT</i> run type.</p> <ul style="list-style-type: none"> • If set to 1, a source analysis only will be performed - the <i>HYPACT</i> run is stopped before the first timestep proceeds with advection and diffusion. The initial output files will contain the initial emission field, providing a useful way for the user to check their work. Note that a summary of each source is also printed to standard output prior to the first timestep. • If set to 2, a normal dispersion run is made • If set to 3, the run will be a backward trajectory run. Only Lagrangian sources are allowed and it is recommended that ihurb be set to 0.
<p>metpref</p> <p>character</p>	<p>The path name and prefix of the names of <i>RAMS</i> analysis files to be read and used to drive the <i>HYPACT</i> simulation (up to 80 characters). Note that the path name is optional and can be either relative or absolute (no path is equivalent to './'). The analysis files to be used must all exist in this same directory, and must span a range of simulation time that brackets the range of time over which a <i>HYPACT</i> simulation is to be run.</p>

<p>dtpart</p> <p>real</p>	<p>The length of the discrete timestep in seconds used to update particle velocities and positions. Its value should be carefully chosen based on required speed and accuracy of the <i>HYPACT</i> run. It would be desirable for each particle to move no more than the dimension of a grid cell (on which the atmospheric data are represented) in a single timestep so that the maximum amount of detail contained in the atmospheric data is utilized. This may at times require a very small value of dtpart, such that a large number of individual timesteps must be performed to run <i>HYPACT</i> for a given span of simulation time. If this runs too slowly, and/or if a lower level of accuracy is acceptable, dtpart may be longer. An estimate for an appropriate value of dtpart can be obtained by dividing grid cell dimensions by atmospheric wind velocity components.</p>
<p>maxpart</p> <p>integer</p>	<p>The total number of particles that can be used in a <i>HYPACT</i> run. An appropriate number can vary widely from one application to another. If one only wants a general idea of which direction particles are traveling from a source, fewer than 1000 particles may be totally adequate. If concentrations are to be computed from the particle distribution, it is desirable to have as many particles as practical to improve the accuracy of the computation. Tens of thousands of particles might be used in this case, or even more if concentrations far from a dense particle plume are sought. Of course, <i>HYPACT</i> runs slower when many particles are activated. Note that maxpart does not specify the number of particles actually emitted (this is done by numparts, described in the \$EMISSIONS namelist) but is the maximum number allowable from computer memory considerations.</p>
<p>freqavg</p> <p>real</p>	<p><i>No longer used. HYPACT will always interpolate in time between the closest existing analysis files in defining the fields for each of the RAMS nested grids. This also takes into account the RAMS v5 feature of allowing different time frequencies of the analysis file output times for different nested grids.</i></p> <p>The time interval in seconds between successive updates of the time average applied to input meteorological data to define it at times close to the current simulation time of the particles in <i>HYPACT</i>. The meteorological data (for example, the output analysis files from <i>RAMS</i>) is defined for discrete points in time, spaced usually at regular intervals such as an hour. Interpolation between consecutive times of the data is performed so that particles use winds and turbulence parameters close to the correct time. Since frequent averaging takes time and is unnecessary freqavg is usually set to a value many times longer than the <i>HYPACT</i> timestep dtpart, but several times shorter than the interval between input data files. Must be less than or equal to the <i>RAMS</i> analysis file frequency, and greater than or equal to dtpart.</p>
<p>ihrturb</p> <p>integer</p>	<p>A flag specifying whether a turbulent component is to be added to the mean resolved atmospheric motion for advecting particles.</p> <ul style="list-style-type: none"> • If set the 0, no turbulence will be included, and the <i>HYPACT</i> Lagrangian particle model operates purely as a trajectory model. • If set to 1, atmospheric turbulent intensity is evaluated and used to derive a turbulent component of motion for each particle. This allows the effects of unresolved atmospheric motions on particle transport and dispersion to be statistically accounted for.

iadvord integer	The order of the advection operator to be used in transporting Eulerian concentration fields. The choices are 2 for second order and 6 for sixth order.
hybfreq real	When an emission type (type) is specified as 'hyb', the Lagrangian particles emitted for the source are converted to Eulerian concentrations at some point downwind (when the computed dispersal of particles is sufficient). This conversion is handled by releasing the particles in clusters. The spread of each cluster is then checked every hybfreq seconds, and if the following criteria is satisfied, the particles are converted to Eulerian concentrations. hybfreq should be set to within a few times dtpart . If hybfreq is less than dtpart , clusters will be checked every <i>HYPACT</i> timestep. The criterion is a measure of spread between the particles compared to the specified horizontal and vertical distances. The spread is taken as six times the standard deviation of cluster particle locations done in each of x, y and z directions. This is then checked against hybhoriz (x and y) and hybvert (z), and the conversion made only if the spread in all three dimensions is greater than the hybhoriz and hybvert criteria.
ihybpert integer	The number of particles emitted per cluster. Different clusters will be allocated to each source so that no cluster will contain a variety of species. ihybpert should be set that so that a cluster of particles does not extend over more than several <i>HYPACT</i> timesteps.
ihybmin integer	The minimum particles per cluster. If there are less than ihybmin particles in a cluster, the cluster will be converted to Eulerian concentrations regardless of the spread. This value must be less than ihybpert , otherwise all particles will be converted when their respective clusters are full. Conversion of a cluster that a source is in the process of filling will not take place until that cluster is full, or the source's emission is complete. A recommended setting for ihybmin is 10% of ihybpert .
hybhoriz real	The horizontal length scale defined as fractional horizontal grid spacing. With a spread of six standard deviations and hybhoriz and hybvert is equal to 1, cluster conversions are made approximately when the particles have dispersed over at least the entire grid volume in each dimension. Settings of hybhoriz and hybvert should therefore be in the range 1 to 3.
hybvert real	The vertical length scale defined as fractional vertical grid spacing or the distance in meters if hybvert is less than 0. With a spread of six standard deviations and hybhoriz and hybvert is equal to 1, cluster conversions are made approximately when the particles have dispersed over at least the entire grid volume in each dimension. Settings of hybhoriz and hybvert should therefore be in the range 1 to 3.

\$OUTPUT Namelist

Almost all output is controlled by the **\$OUTPUT** namelist, with the sole exception that the [ioutspec](#) variable in the [\\$SPECIES namelist](#) controls whether individual species are included in the output.

hyppref character	The UNIX path name and prefix for the HYPACT output files generated during the simulation (up to 80 characters). Note that the path is optional and can be either relative or absolute (no path is equivalent to './'). The time interval between the writing of successive files in the set is specified by the namelist variable hypfreq .
hypfreq real	The time interval in seconds between successive writing of HYPACT output files. The settings of ipartout , ieulout and ilagout specify what variables are output, and irgrid specifies the RAMS grids that are included in the output. Output file names and formats are described separately following the namelist documentation.
ipartout integer	Particle file output flag. ipartout set to 1 produces particle files every hypfreq seconds provided at least one element of type is equal to 'lag', 'hyb' or 'both'.
ieulout ilagout integer	Output concentrations are controlled with the following, noting that there is one concentration field for each species; <ul style="list-style-type: none"> • ieulout writes the Eulerian concentrations. These fields will also contain the concentrations produced by the conversion of particle clusters in hybrid mode (\$EMISSIONS namelist parameter type set to 'hyb'). • ilagout writes the concentration fields computed by converting all particles to Eulerian concentrations (the particles still remain in the Lagrangian field). For each: <ul style="list-style-type: none"> • 0 gives no output. • 1 gives the concentrations at the end the current timestep. • 2 gives the average concentrations over avgtime seconds prior to the file output time (determined by hypfreq). • 3 gives both the current and average concentration output. The hybrid concentrations (when an emission type (type) is specified as 'hyb') are the sum of the Lagrangian particles yet to be converted and Eulerian concentrations from converted clusters for each species. This summation is handled in the post processing with REVU . All concentrations are computed and output on the input RAMS grid structure.
imetout integer	Meteorology output control flag. imetout set to 1 outputs the interpolated RAMS meteorology as used by HYPACT at the output time. The meteorology is included in the concentration file output.

avgtime real	Averaging time in seconds over which concentration field averaging is done before each output time. Averaging can be done for any species on the <i>RAMS</i> grid. It cannot be done for individual emission sources unless each emission source emits a different species, and cannot be done for the particle files. avgtime must be less than or equal to the \$GENERAL namelist parameter hypfreq .
----------------------------	--

ioutfmt integer	Output file format: <ul style="list-style-type: none">• If set to 1, VFILE binary files are output.
irgrid integer	<p>HYPACT output is limited to the grids specified by irgrid.</p> <ul style="list-style-type: none">• If set to 0, concentrations and particles are output on all grids.• If set to greater than 0, concentrations are output for that grid only.• If set to less than 0, concentrations are output for all grids less than the absolute value of irgrid. <p>In the latter two cases, particles in the output file include all those that reside within the bounds of the specified irgrid.</p>

\$SPECIES Namelist

specfile character	The path and filename containing the species database (up to 80 characters). An example file and the file format follows this table. Note that the path is optional and can be either relative or absolute (no path is equivalent to './'). If specfile is set to 'none', the species database is read from the following namelist parameters. If specfile points to a database file, with the exception of ioutspec , the remaining \$SPECIES namelist parameters are overwritten.
specname character array [†]	A multi-valued parameter specifying the name of each species (up to 30 characters long). specname can include standard NCAR Graphics sub and superscripting code. This information is stored in the output 'SPEC' file, and so it is useful for later plotting with NCAR Graphics. 'N' denotes the beginning of normal text, 'B' subscripted text and 'S' superscripted text (all ":" delimited). It is safest to start and end with an 'N', for example, ':N:CO:B:2:N:' results in CO ₂ .
wgtmol real array [†]	A multi-valued parameter specifying the molecular weight of each species in grams per mole. wgtmol is used to calculate Eulerian concentrations in parts per trillion if units is in nanograms (ng), parts per million (ppm) if units is in micrograms (ug) and parts per thousand if units is in milligrams (mg).
units character array [†]	A multi-valued parameter specifying the mass units applied to the emission rate and ratio (up to 30 characters long). units can be set to ng (nanograms), ug (micrograms) or mg (milligrams). The units set here should be consistent with those used for the \$EMISSIONS namelist parameters, rate and ratio .
ihfall integer array [†]	A flag specifying whether particles in <i>HYPACT</i> are to be caused by gravity to settle relative to the local flow in which they are embedded. A value of 0 indicates no settling, while a value of 1 activates the settling. The parameters szmin , szmax , and szpwr below control the settling velocity for particles from each source.
szmin real array [†]	A multi-valued parameter specifying a minimum particle diameter in meters emitted from each source region. Its only use is in determining gravitational settling speed, and is ignored if namelist variable ihfall is set to 0.
szmax real array [†]	A multi-valued parameter corresponding to namelist variable szmin , but applying to the maximum particle diameter from each source region. Currently, only a Stoke's drag formula is implemented in <i>HYPACT</i> , which is relevant for diameters up to about than 80×10^{-6} m.
szpwr real array [†]	A multi-valued parameter relating closely to szmin and szmax , and is used in computing gravitational settling speed. It specifies a particle size distribution, effectively weighting the particles toward the larger or smaller sizes. szpwr specifies the power of the diameter which is linearly (uniformly) distributed between the specified size limits. For example, if szpwr is set to 1, all diameters between the limits are emitted with equal probability. If szpwr is set to 3, all cubes (third power) of particle diameters, or equivalently all volumes or masses, are emitted with equal probability between the size limits. szpwr may be any non-negative number.

ioutspec integer array [†]	A multi-valued parameter specifying what species are to be included in the output. To output the species, set ioutspec to 1.
---	---

[†] Each array must contain entries for each species.

Species Database File

The following illustrates some species in the database file format:

```
1, 3
TRS, 60, ug, 0, 1., 1., 1.
:N:CO:B:2:N:, 44, ug, 0, 1., 1., 1.
XXT, 98, ug, 0, 1., 1., 1.
```

The header line contains file format version and the number of species in the file with each subsequent line containing all the information for each species in comma delimited format (order is important):

- [specname](#), [wgtmol](#), [units](#), [ihfall](#), [szpwr](#), [szmin](#), [szmax](#)

Note that the [\\$SPECIES namelist](#) variable [ioutspec](#) is not included in the species database file, and should be included in the [\\$SPECIES namelist](#). If it is not, the default is to output for all species.

\$\$SOURCES Namelist

srcfile character	The path and filename containing the source database (up to 80 characters). An example file and the file format follows this table. Note that the path is optional and can be either relative or absolute (no path is equivalent to './'). If srcfile is set to 'none' the source database is read from the following namelist parameters. If srcfile points to a database file the remaining \$\$SOURCES namelist parameters are overwritten.
srcname character array [†]	A multi-valued parameter specifying the name of each source (up to 30 characters long).
shape character array [†]	<p>A multi-valued parameter specifying the horizontal cross-sectional shape of each source. shape can be set to 'point', 'rectangle', 'triangle', 'polygon', or 'ellipse' (these can be abbreviated to the first three characters).</p> <ul style="list-style-type: none"> • If a point source is specified, namelist variables srcx and srcy are used to determine the horizontal location of the source, and srcz the height of the source (this is converted to a rectangular source with no dimensions in <i>HYPACT</i>). • If a triangular or irregular polygon shape is specified, the namelist variables polylat and polylon are used to specify the horizontal coordinates of the vertices, srcx, srcy the center of the polygon (see the notes on polylat and polylon for the relevance of this), and srcz and zsize the height and thickness of the source. • If a rectangular or elliptical shape is specified, namelist variables srcx, srcy, xsize, and ysize are used to determine the horizontal location and size of the source, and srcz and zsize the height and thickness of the source. rotation can also be specified to rotate rectangular or elliptical sources in the horizontal plane. • If a gridded shape is specified, the emissions are obtained from the file named in gridfile. (not implemented) <p>Note that the center location of each source, and the entire source region, should lie within the volume over which the input meteorological fields are defined (i.e., the coarsest grid of the <i>RAMS</i> domain). Any location outside this range places the source outside the model domain, and emissions there will remain stationary in <i>HYPACT</i>.</p>
srcx srcy real array [†]	Multi-valued parameters specifying the horizontal coordinates of the center of each source region. srcx and srcy are used only for point, rectangular, polygonal and elliptical sources specified by shape , and indicate the geographic longitude and latitude, respectively, in degrees.
srcz real array [†]	A multi-valued parameter specifying the height above the ground in meters, and applies for all source types.

<p>polylat polylon</p> <p>real array[†]</p>	<p>Multi-valued two-dimensional parameters specifying the geographic vertex latitude and longitude coordinates for triangular and irregular polygon shaped sources specified by shape. The indices on the 2-D array are vertex and source. The vertex index must always be equal to 1 (first array element), while the source index refers to the source number. For irregular polygon sources, the vertices must be defined in an anti-clockwise order - the first and last coordinates are joined to enclose the polygon. For example, the following describes a triangular source for source 2, and a four-sided polygon for source 3:</p> <pre>polylat(1,2) = 40.075, 40.076, 40.071, polylon(1,2) = -104.721, -104.723, -104.722, polylat(1,3) = 40.075, 40.075, 40.071, 40.071, polylon(1,3) = -104.721, -104.723, -104.723, -104.721,</pre> <p>Polygonal sources were conceived to facilitate simpler specification of irregularly shaped sources, such as urban emission boundaries. Polygonal sources are broken up into an array of triangular sources in the <i>HYPACT</i> code. Each defined polygon must be sufficiently simple so that none of the triangles overlap, and no part of a triangle lies outside the polygon. Tests for these limitations are made in the code. In breaking up a polygon, the vertices of each triangle are initially taken as the two points of one side of the polygon, and a point somewhere in the center of the polygon (as defined by srcx and srcy). If any triangle fails a test, the central point is moved in the code, and the triangles redefined. If the code generated triangles continue to fail a test, the program is stopped and the user is required to reassess the configuration of their sources.</p>
<p>xsize ysize</p> <p>real array[†]</p>	<p>Multi-valued parameters specifying the x and y direction dimensions in meters of rectangular and elliptical sources specified by shape.</p> <ul style="list-style-type: none"> • If rectangular, xsize and ysize are the dimensions respectively. • If elliptical, xsize and ysize represent the lengths of the major and minor axes respectively.
<p>zsize</p> <p>real array[†]</p>	<p>Similar to xsize and ysize, but specifying the vertical dimension of the source in meters. If zsize is greater than zero, half of the source will be below the specified source center height srcz. Thus, the latter should be sufficiently high that no part of the source is underground (i.e., at least half the value of srcz).</p>
<p>rotation</p> <p>real array[†]</p>	<p>A multi-valued variable indicating a rotation angle (in degrees) in a clockwise manner about a vertical axis of a source. This allows basic shapes such as rectangle or ellipse to be first indicated with principal axes parallel to the x- and y-axes of the <i>HYPACT</i> grid, and then for those shapes to be re-oriented in any direction desired. A positive angle causes a clockwise rotation as viewed from above.</p>

[†] Each array must contain entries for each source, noting the [flexibility demonstrated in the example](#) below.

Source Namelist Examples

The following gives an example for each type of source. Note that different parameters apply in different ways to the sources, and some placeholders may be required, although they will not get used. There are two recommended ways to specify sources - either by using placeholders:

```

srcname='Mill_1', 'Power_1', 'Power_2', 'Ponds', 'Mill_2',
shape='rec', 'tri', 'poly', 'ell', 'poi',
srcy=40.2, 0., 39.7, 40., 40.2,
srcx=-104.2, 0., -82.7, -105., -105.2,
srcz=50., 100., 100., 5., 50.,
xsize=100., 0., 0., 8000., 0.,
ysize=100., 0., 0., 8000., 0.,
zsize=15., 10., 100., 10., 0.,
polylat(1,2)=40.075, 40.076, 40.071,
polylon(1,2)=-104.721, -104.723, -104.722,
polylat(1,3)=40.075, 40.075, 40.071, 40.071,
polylon(1,3)=-104.721, -104.723, -104.724, -104.720,
rotation=45., 0., 0., 40., 0.,

```

Note the placeholders:

- The second values of srcy and srcx are not used for the triangular source.
- The second, third and fifth values of xsize, ysize and rotation are not used for the triangular, polygonal and point sources.

Note also that only the second and third series of polylat and polylon points (for the triangular and polygonal sources) are required. In this case the two dimensional namelist variables must contain the relevant indices (the second index refers to the source number).

Or alternatively, to remove the need for placeholders - each source can be specified with the relevant indices on the namelist variables:

```

srcname(1)='Mill_1',
shape(1)='rec',
srcy(1)=40.2,
srcx(1)=-104.2,
srcz(1)=50.,
xsize(1)=100.,
ysize(1)=100.,
zsize(1)=15.,
rotation(1)=45.,

srcname(2)='Power_1',
shape(2)='tri',
srcz(2)=100.,
polylat(1,2)=40.075, 40.076, 40.071,
polylon(1,2)=-104.721, -104.723, -104.722,
zsize(2)=10.,

```

```

srcname(3)='Power_2',
shape(3)='poly',
srcy(3)=39.7,
srcx(3)=-82.7,
srcz(3)=100.,
polylat(1,3)=40.075,40.075,40.071,40.071,
polylon(1,3)=-104.721,-104.723,-104.724,-104.720,
zsize(3)=100.,

srcname(4)='Ponds',
shape(4)='ell',
srcy(4)=40.,
srcx(4)=-105.,
srcz(4)=5.,
xsize(4)=8000.,
ysize(4)=8000.,
zsize(4)=0.,
rotation(4)=40.,

srcname(5)='Mill_2',
shape(5)='poi',
srcy(5)=40.2,
srcx(5)=-105.2,
srcz(5)=5.,

```

Source Database File

The following illustrates the above sources in the database file format:

```

1,5
Mill_1,rec,40.,-83,50.,1.,1.,1.,0.
Power_1,tri,40.,-83,100.,40.075,-104.721,40.075,-104.723 \
,40.071,-104.722,0.
Power_2,poly,40.,-83,100.,4,40.075,-104.721,40.075,-104.723 \
,40.071,-104.724,40.071,-104.720,0.
Ponds,ell,40.,-83,500.,300000.,300000.,100.,0.
Mill_2,poi,40.,-83,5.

```

Where the header line contains file format version and the number of sources in the file and one line for each source containing all the information for each source in comma delimited format. The relevant information for each source shape is (order is important):

- Rectangular: [srcname](#), [shape](#), [srcy](#), [srcx](#), [srcz](#), [xsize](#), [ysize](#), [zsize](#), [rotation](#)
- Triangular: [srcname](#), [shape](#), [srcz](#), [polylat](#), [polylon](#), [zsize](#)
- Polygonal: [srcname](#), [shape](#), [srcy](#), [srcx](#), [srcz](#), [nvert](#), [polylat](#), [polylon](#), [zsize](#)
- Elliptical: [srcname](#), [shape](#), [srcy](#), [srcx](#), [srcz](#), [xsize](#), [ysize](#), [zsize](#), [rotation](#)
- Point: [srcname](#), [shape](#), [srcy](#), [srcx](#), [srcz](#)

Note the extra variable **nvert** in the polygonal source that is not in the namelist. This is the number of vertices in the polygon.

\$EMISSIONS Namelist

emfile character	The path and filename containing the emission database (up to 80 characters). An example file and the file format follows this table. Note that the path is optional and can be either relative or absolute (no path is equivalent to './'). If emfile is set to 'none', the emission database is read from the following namelist parameters. If specfile points to a database file, with the exception of iemit and csimend , the remaining \$EMISSIONS namelist parameters are overwritten.
iemit integer array [†]	A multi-valued parameter that determines what emissions specified by this namelist are to be used for the current HYPACT simulation. A value of 1 activates the emission.
isource integer array [†]	A multi-valued parameter that specifies the nth source from the \$SOURCES namelist or source file .
ispecies integer array [†]	A multi-valued parameter that specifies the nth species from the \$SPECIES namelist or species file .
crelstrt character array [†]	<p>A multi-valued parameter which specify the UTC date/time when the emissions for each source begins.</p> <p>The format of the character string is: yyyy-mm-dd-hh:mn:ss to specify the year, month, day, hour, minutes, seconds. An example would be: '2000-07-30-18:00:00', specifying 30 July 2000 at 1800.00 UTC.</p> <p>The earliest time specified by irelstrt for an emission that is activated by iemit defines the start time of the HYPACT simulation.</p>
crelend character array [†]	A multi-valued parameters that specify the ending time of particle and/or concentration emission from each source. It is related closely to irelstrt . creldur uses the same syntax for the number of hours, minutes, and seconds as irelstrt . Together, irelstrt and crelend determine the duration for the emission of each source. A duration time of zero or anything less than the timestep length dtpart causes a source to act as in instantaneous or “puff” release.
csimend character	Parameter that specifies the ending time of the HYPACT simulation. csimend is formatted as described for namelist parameter crelstrt .

<p>type</p> <p>character array[†]</p>	<p>A multi-valued parameter that specifies the source type.</p> <ul style="list-style-type: none">• If set to 'lag' (Lagrangian), only Lagrangian particles will be released from the source.• If set to 'eul' (Eulerian), only Eulerian concentrations will be released from the source.• If set to 'hyb' (hybrid), Lagrangian particles will be released, then transformed into Eulerian concentrations when the hybrid criteria (ihymin, hybhoriz and hybvert) set in the \$GENERAL namelist are met.• If set to 'both' (both), both Eulerian concentrations and Lagrangian particles will be released from the source. Fields will be calculated as if there were two sources (i.e., concentrations will be effectively doubled). This is primarily used for testing purposes.
---	---

<p>scaling</p> <p>character array[†]</p>	<p>A multi-valued parameter that specifies how the emission rate and number of particles are determined. Any two of the three namelist variables rate, numparts and ratio can be used to calculate the third as follows:</p> <ol style="list-style-type: none"> If set to 'rat' (rate), the <u>emission rates are determined</u> from the particle release rate (numparts), the size of the source (volume, area or length determined from xsize, ysize, zsize and shape from the \$SOURCES namelist - noting that the size is set to 1 for a point source), the timestep (dtpart from the \$GENERAL namelist), and the mass per particle ratio (ratio): $\text{emission rate} = \text{particle rate} * \frac{\text{mass ratio}}{\text{size} * \text{timestep}}$ <p>The particle rate (numparts) has units of particles/timestep. If the mass units are ug (units from the \$SPECIES namelist) the mass ratio (ratio) has units of ug/particle. Thus, if the source is volumetric (size has units of m³), the emission rate has units of (ug/s)/m³ (i.e., it is a specific rate that is calculated).</p> If set to 'tot' (total), numparts is the total number of particles released and the particle release rate is determined from the total particles, the release duration (determined from ireldur, and idurdays) and the timestep: $\text{particle rate} = \text{total particles} * \frac{\text{timestep}}{\text{release duration}}$ <p>The <u>emission rates are then determined</u> as for scaling set to 'rat'.</p> If set to 'spe' (specific), the <u>particle release rates are determined</u> as a function of the specific emission rate (rate), the source size, the timestep and the mass ratio: $\text{particle rate} = \text{emission rate} * \text{size} * \frac{\text{timestep}}{\text{mass ratio}}$ <p>The particle rate has units of particles/s.</p> If set to 'abs' (absolute), the <u>particle release rates are determined</u> from absolute emission rate (rate), the timestep and the mass ratio: $\text{particle rate} = \text{emission rate} * \text{timestep} / \text{mass ratio}$ <p>The calculate particle rate has units of particles/s.</p> If set to 'spe' (specific) and the mass ratio (ratio) is set to 0, the <u>mass ratios are determined</u> as a function of the specific emission rate (rate), the size of the source, the timestep and the particle release rates (numparts): $\text{mass ratio} = \text{emission rate} * \text{size} * \frac{\text{timestep}}{\text{particle rate}}$ <p>If the mass units in the emission rate are defined as 'ug', the derived mass ratio has units of ug/particle.</p> If set to 'abs' (absolute) and the mass ratio (ratio) is set to 0, the <u>mass ratios are determined</u> from absolute emission rate (rate), the timestep and the particle release rates (numparts): $\text{mass ratio} = \text{emission rate} * \frac{\text{timestep}}{\text{particle rate}}$ <p>If the mass units in the emission rate are defined as 'ug', the derived mass ratio has units of ug/particle.</p> <p>Note that, as whole numbers of particles only can be emitted, particle emission rates from timestep to timestep are adjusted to maintain correct average emission rates. A set of emissions examples is included below.</p>
--	--

rate real array [†]	<p>A multi-valued parameter that specifies the emission rates if scaling is equal to 'spe' or 'abs'.</p> <ul style="list-style-type: none"> If scaling is equal to 'spe', rate is the specific emissions rate. In this case, if the mass units in the emission rate are defined as 'ug', and the source is volumetric, the units of rate are (ug/s)/m³. If scaling is equal to 'abs', rate is the absolute emissions rate. In this case, if the mass units in the emission rate are defined as 'ug', the units of rate are ug/s. <p>If scaling is equal to 'rat', the emission rate is calculated. The same mass units must be used for rate and ratio, which should also be consistent with the \$SPECIES namelist variable, units.</p>
numparts integer array [†]	<p>A multi-valued parameter that specifies the number of particles emitted per <i>HYPACT</i> timestep (dtpart) if scaling is equal to 'rat', or if scaling is equal to 'spe' or 'abs' and ratio is equal to 0. If scaling is equal to 'spe' or 'abs' and ratio has a positive value, the number of particles is calculated. Note that the total particles emitted for all sources cannot exceed the \$GENERAL namelist variable, maxpart.</p>
ratio real array [†]	<p>A multi-valued parameter which specifies the mass to particle ratio used to calculate the number of particles released per timestep if scaling is equal to 'sca', or the emission rate if scaling is equal to 'tot'. If ratio is set 0 the mass ratio is calculated. The same mass units must be used for ratio and rate, which should also be consistent with the \$SPECIES namelist variable, units.</p>

[†] Each array must contain entries for each emission.

Emissions Database File

The following illustrates some emissions in the database file format:

1,4

```
4,2,120000,0,060000,0,both,rat,20,20.,1.e9
1,2,120000,0,060000,0,lag,abs,1,.05,120.
1,2,130000,0,020000,0,lag,rat,5,1.,0.
4,1,130000,0,020000,0,lag,rat,3,20.,1.e9
```

Where the header line contains file format version and the number of emissions in the file with each subsequent line containing all the information for each emission in comma delimited format (order is important):

- [isource](#), [ispecies](#), [irelstrt](#), [istrtdays](#), [ireldur](#), [idurdays](#), [type](#), [scaling](#), [numparts](#), [rate](#), [ratio](#)

Note that the [\\$EMISSIONS namelist](#) variables [iemit](#), [isimend](#) end [enddays](#) are not included in the emissions database file, and should always be included in the [\\$EMISSIONS namelist](#). If they are not, the defaults are to emit all emissions and run the *HYPACT* simulation to the end of the *RAMS* run.

Emissions Examples

Preliminary Notes

- If memory and computational speed are issues, you will want to pay close attention to the numbers of particle released in a simulation.
- When converting Lagrangian particles to concentrations you want to ensure that the mass ratio of a particle results in a sufficiently defined concentration field.

For example, if the sensors you are comparing the simulation results with are good to 4 ug/m³ and your grid size at the ground surface is 100 m square and 25 m deep (250,000 m³), then you do not want your particle mass ratio to exceed:

$$\begin{aligned} \text{mass ratio} &= 250,000 \text{ m}^3 * \frac{4 \text{ ug}}{\text{m}^3} * \frac{1}{1 \text{ particle}} \\ &= 1,000,000 \text{ ug} / \text{particle} = 1 \text{ kg} / \text{particle} \end{aligned}$$

This is not as an important factor is determining a Eulerian source as the particles are only used as a convenient method for statistically distributing the emissions throughout the source volume or area (noting that a source can be complex in shape and cross the boundaries of any number of grid cells).

This implies that you should have a large number of particles configured for a Eulerian source as:

1. This will give a statistically even distribution of particles through the source region (and no false peaks).
2. They are converted immediately to concentrations upon emission and do not accumulate in the model memory.

This is not as an important factor is determining a Eulerian source as the particles are only used as a convenient method for statistically distributing the emissions throughout the source volume or area (noting that a source can be complex in shape and cross the boundaries of any number of grid cells).

- When computing the various quantities for each emission it is a good idea to run **HYPACT** in its source analysis mode ([ihyprun](#) = 1) as the full emission summary is computed and printed to screen.

Case Study Scenario

Six 10 by 10 m cooling tower plumes elevated 25 m from the ground emitting at a rate of 0.2 kg/s.

In the examples that follow, each emission specifies the cooling tower with different parameters known and unknown. In all cases, the computed emission parameters compute to the save values for input into **HYPACT**, as illustrated by the source summaries.

Example #1 - Rate

Typical Usage

If you *do not know the emission rate*, as is often the case with arbitrary sources, the 'rate' scaling option is a good way to get a sensible emission with a known amount of particles. In this way you can ensure that you release enough particles per timestep so that you get good concentration definition in the region of most interest from a Lagrangian particle source.

Relevant Namelist Settings

```

dtpart = 50 s
xsize = 10 m
ysize = 10 m
zsize = 0 m
units = ug
duration = 21600 s (6 hours as specified by crelstrt, crelend)
scaling = rat(e)
numparts = 10 particles/timestep
rate = 0 ug/m2/s (UNKNOWN - calculated by HYPACT)
ratio = 1,000,000 ug/particle

```

emission rate = particle rate * mass ratio / (size * timestep)

$$\begin{aligned}
 \text{rate} &= \frac{\text{parts} * \text{ratio}}{\text{size} * \text{timestep}} = \frac{10 \text{ parts}}{\text{timestep}} * \frac{1,000,000 \text{ ug}}{\text{particle}} * \frac{1}{100 \text{ m}^2} * \frac{\text{timestep}}{50 \text{ s}} \\
 &= 2,000 \frac{\text{ug}}{\text{m}^2 \text{ s}} \\
 &= 200,000 \text{ ug} / \text{s} = 0.2 \text{ kg} / \text{s}
 \end{aligned}$$

Source Summary Output

```

Model Emission Source: 1
Source: 1 ISP Species: 1 :N:CO:B:2:N:
Type definition - both rat
Source shape - square
Location(lat/long) - 40.79390 -73.10170
Centre (x,y,z) - 75573.16 33045.85 100.0000 m
Height Dimension - 0.000000 m
Horizontal Dimensions - 10.00000 x 10.00000 m2
Size - 100.0000 m2
Emission times:
start - 0.000000 s
end - 21600.00 s
duration - 21600.00 s
Timestep - 50.00000 s
Timesteps - 432.0000
Particles to release - 10.00000 per timestep
- 4320.000 total for this source
Mass per particle - 1000000. ug/particle
Emission rate - 2000.000 ug/m2/s
- 200000.0 ug/s
- 1.0000000E+07 ug/timestep
- 4.3200000E+09 ug (total for this source)

```

Example #2 - Total

Typical Usage

If you *do not know the emission rate*, as is often the case with arbitrary sources, the 'total' scaling option is a good way to get a sensible emission with a known total amount of particles. In this way you can ensure that you release enough particles per timestep so that you get good concentration definition in the region of most interest from a Lagrangian particle source.

Relevant Namelist Settings

```

dtpart = 50 s
xsize = 10 m
ysize = 10 m
zsize = 0 m
units = ug
duration = 21600 s (6 hours as specified by ireldur, idurdays)
scaling = tot(al)
numparts = 4320 particles (total released fro this emission)
rate = 0 ug/m2/s (UNKNOWN - calculated by HYPACT)
ratio = 1,000,000 ug/particle

```

particle rate = total particles * timestep / release duration

$$\text{particle rate} = 4320 \text{ parts} * \frac{50 \text{ s}}{\text{timestep}} * \frac{1}{21600 \text{ s}} = 10 \frac{\text{parts}}{\text{timestep}}$$

emission rate = particle rate * mass ratio / (size * timestep)

$$\begin{aligned} \text{rate} &= \frac{\text{parts} * \text{ratio}}{\text{size} * \text{timestep}} = \frac{10 \text{ parts}}{\text{timestep}} * \frac{1,000,000 \text{ ug}}{\text{particle}} * \frac{1}{100 \text{ m}^2} * \frac{\text{timestep}}{50 \text{ s}} \\ &= 2,000 \text{ ug} / \text{m}^2 \text{ s} = 200,000 \text{ ug} / \text{s} = 0.2 \text{ kg} / \text{s} \end{aligned}$$

Source Summary Output

```

Model Emission Source: 2
Source: 2 MTP Species: 1 :N:CO:B:2:N:
Type definition - both tot
Source shape - square
Location(lat/long) - 41.07310 -71.92330
Centre (x,y,z) - 173972.5 65746.22 25.00000 m
Height Dimension - 0.000000 m
Horizontal Dimensions - 10.00000 x 10.00000 m2
Size - 100.0000 m2
Emission times:
  start - 0.000000 s
  end - 21600.00 s
  duration - 21600.00 s
Timestep - 50.00000 s
Timesteps - 432.0000
Particles to release - 10.00000 per timestep
- 4320.000 total for this source
Mass per particle - 1000000. ug/particle
Emission rate - 2000.000 ug/m2/s
- 200000.0 ug/s
- 1.0000000E+07 ug/timestep

```

- 4.3200000E+09 ug (total for this source)

Example #3 - Specific

Typical Usage

Used when you have a *maximum specification for the particle mass ratio* (see the [preliminary note](#) above), and particularly useful when you know the release rate per m² (or in whatever spatial units the source is described in). However, you do not have or require control on the number of particles being released.

Relevant Namelist Settings

```

dtpart = 50 s
xsize = 10 m
ysize = 10 m
zsize = 0 m
units = ug
duration = 21600 s (6 hours as specified by ireldur, idurdays)
scaling = spe(cific)
numparts = 0 particles/s (UNKNOWN - calculated by HYPACT)
rate = 2,000 ug/m2/s
ratio = 1,000,000 ug/particle

```

```
particle rate = emission rate * size * timestep / mass ratio
```

$$\begin{aligned}
 \text{particle rate} &= \frac{2,000 \text{ ug}}{\text{m}^2 \text{ s}} * 100 \text{ m}^2 * \frac{50 \text{ s}}{\text{timestep}} * \frac{\text{particle}}{1,000,000 \text{ ug}} \\
 &= 10 \text{ parts} / \text{timestep}
 \end{aligned}$$

Source Summary Output

```

Model Emission Source: 3
Source: 3 LGA Species: 1 :N:CO:B:2:N:
Type definition - both spe
Source shape - square
Location(lat/long) - 40.77920 -73.88000
Centre (x,y,z) - 10097.79 31032.92 25.00000 m
Height Dimension - 0.000000 m
Horizontal Dimensions - 10.00000 x 10.00000 m2
Size - 100.0000 m2
Emission times:
  start - 0.000000 s
  end - 21600.00 s
  duration - 21600.00 s
Timestep - 50.00000 s
Timesteps - 432.0000
Particles to release - 10.00000 per timestep
- 4320.000 total for this source
Mass per particle - 1000000. ug/particle
Emission rate - 2000.000 ug/m2/s
- 200000.0 ug/s
- 1.0000000E+07 ug/timestep
- 4.3200000E+09 ug (total for this source)

```

Example #4 - Absolute

Typical Usage

Used when you have a *maximum specification for the particle mass ratio* (see the [preliminary note](#) above), and particularly useful with complex source shapes where you know the total release rate. However, you do not have or require control on the number of particles being released.

Relevant Namelist Settings

```

dtpart = 50 s
xsize  = 10 m
ysize  = 10 m
zsize  = 0 m
units  = ug
duration = 21600 s (6 hours as specified by ireldur, idurdays)
scaling = abs(olute)
numparts = 0 particles/s (UNKNOWN - calculated by HYPACT)
rate     = 200,000 ug/s
ratio    = 1,000,000 ug/particle

```

```
particle rate = emission rate * timestep / mass ratio
```

$$\begin{aligned}
 \text{particle rate} &= \frac{200,000 \text{ ug}}{\text{s}} * \frac{50 \text{ s}}{\text{timestep}} * \frac{\text{particle}}{1,000,000 \text{ ug}} \\
 &= 10 \text{ parts} / \text{timestep}
 \end{aligned}$$

Source Summary Output

```

Model Emission Source: 4
Source: 4 POU Species: 2 :N:S0:B:2:N:
Type definition - both abs
Source shape - square
Location(lat/long) - 41.62560 -73.88190
Centre (x,y,z) - 9810.993 125092.3 25.00000 m
Height Dimension - 0.000000 m
Horizontal Dimensions - 10.00000 x 10.00000 m2
Size - 100.0000 m2
Emission times:
start - 0.000000 s
end - 21600.00 s
duration - 21600.00 s
Timestep - 50.00000 s
Timesteps - 432.0000
Particles to release - 10.00000 per timestep
- 4320.000 total for this source
Mass per particle - 1000000. ug/particle
Emission - 200000.0 ug/s
- 2000.000 ug/m2/s
- 1.0000000E+07 ug/timestep
- 4.3200000E+09 ug (total for this source)

```

Example #5 - Specific

Typical Usage

Used when the *particle mass ratio is unknown* and particularly useful when you know the release rate per m² (or in whatever spatial units the source is described in). You also have control on the number of particles being released per timestep.

Relevant Namelist Settings

```

dtpart = 50 s
xsize = 10 m
ysize = 10 m
zsize = 0 m
units = ug
duration = 21600 s (6 hours as specified by ireldur, idurdays)
scaling = spe(cific)
numparts = 10 particles/s
rate = 2,000 ug/m2/s
ratio = 0 ug/particle (UNKNOWN - calculated by HYPACT)

```

mass ratio = emission rate * size * timestep / particle rate

$$\begin{aligned}
 \text{mass ratio} &= \frac{2,000 \text{ ug}}{\text{m}^2 \text{ s}} * 100 \text{ m}^2 * \frac{50 \text{ s}}{\text{timestep}} * \frac{\text{s}}{10 \text{ particles}} \\
 &= 1,000,000 \text{ ug / particle} = 1 \text{ kg / particle}
 \end{aligned}$$

Source Summary Output

```

Model Emission Source: 5
Source: 5 BLM Species: 2 :N:S0:B:2:N:
Type definition - both spe
Source shape - square
Location(lat/long) - 40.18330 -74.13330
Centre (x,y,z) - -11317.39 -35184.75 25.00000 m
Height Dimension - 0.000000 m
Horizontal Dimensions - 10.000000 x 10.00000 m2
Size - 100.0000 m2
Emission times:
start - 0.000000 s
end - 21600.00 s
duration - 21600.00 s
Timestep - 50.000000 s
Timesteps - 432.0000
Particles to release - 10.00000 per timestep
- 4320.000 total for this source
Mass per particle - 1000000. ug/particle
Emission rate - 2000.000 ug/m2/s
- 200000.0 ug/s
- 1.0000000E+07 ug/timestep
- 4.3200000E+09 ug (total for this source)

```

Example #6 - Absolute

Typical Usage

Used when the *particle mass ratio is unknown* and particularly useful with complex source shapes where you know the total release rate. You also have control on the number of particles being released per timestep.

Relevant Namelist Settings

```

dtpart = 50 s
xsize = 10 m
ysize = 10 m
zsize = 0 m
units = ug
duration = 21600 s (6 hours as specified by ireldur, idurdays)
scaling = abs(olute)
numparts = 10 particles/s
rate = 200,000 ug/s
ratio = 0 ug/particle (UNKNOWN - calculated by HYPACT)

```

mass ratio = emission rate * timestep / particle rate

$$\text{mass ratio} = \frac{200,000 \text{ ug}}{\text{s}} * \frac{50 \text{ s}}{\text{timestep}} * \frac{\text{timestep}}{10 \text{ particles}}$$

$$= 1,000,000 \text{ ug / particle} = 1 \text{ kg / particle}$$

Source Summary Output

```

Model Emission Source: 6
Source: 6 ACY Species: 2 :N:S0:B:2:N:
Type definition - both abs
Source shape - square
Location(lat/long) - 39.46470 -74.58670
Centre (x,y,z) - -50337.31 -114885.1 25.00000 m
Height Dimension - 0.000000 m
Horizontal Dimensions - 10.00000 x 10.00000 m2
Size - 100.0000 m2
Emission times:
start - 0.000000 s
end - 21600.00 s
duration - 21600.00 s
Timestep - 50.00000 s
Timesteps - 432.0000
Particles to release - 10.00000 per timestep
- 4320.000 total for this source
Mass per particle - 1000000. ug/particle
Emission - 200000.0 ug/s
- 2000.000 ug/m2/s
- 1.0000000E+07 ug/timestep
- 4.3200000E+09 ug (total for this source)

```


HYPACT Output Files

Filename convention notes:

1. "*" following filenames represents the date string yyyy-mm-dd-hhmm.
2. "#" represents the grid number.
3. In the following examples the [\\$OUTPUT namelist](#) variable [hyppref](#) = 'a'.

HYPACT Specification File

FORTRAN free format ASCII file *a-spec.txt*.

Main header line:

```
File Version ..... i
Number of Sources ..... npsources
Number of Species ..... nspecies
Average Source Location (lat, lon) .. xavg, yavg
Average Source Height ASL ..... zavg
HYPACT Run Start ..... iy, imn, id, ih, im
```

Then for each source:

```
Source ID ..... psource(nsrc)
Source Type ..... type(nsrc)
Source Name ..... srcname(nsrc)(1:indx1)
Source Location (lat, lon) ..... srcy(nsrc), srcx(nsrc)
Source Height (AGL, ASL) ..... srcz(nsrc), zasl(nsrc)
Species ID ..... pspecies(species(nsrc))
Species Name ..... specname(species(nsrc))
Mass Units ..... units(species(nsrc))
```

Hybrid Grid File

FORTRAN free format ASCII file *a-hgrid.txt*.

Main header line:

```
File Version ..... i
Grid Dimensions (points in x, y, z) .. nih, njh, nk
Grid Location (lat, lon) ..... pslat, pslon
```

Followed by four blocks:

```
x Grid Coordinates (m) ..... xh(i) (i = 1 to nih)
y Grid Coordinates (m) ..... yh(j) (j = 1 to njh)
Grid Heights ASL (m) ..... zh(j) (k = 1 to nk)
Topography Heights (m) ..... topth(i,j) (i = 1 to nh1, j = 1 to nh2)
Land Percentages ..... pctlh(i,j) (i = 1 to nh1, j = 1 to nh2)
```

HYPACT Particle Files

FORTRAN formatted ASCII Files (*a-P-**)

Main header line (format *i3,e16.8,8i8*):

```
File Version ..... i
HYPACT Time (s) ..... hytime
Year ..... iy
Month ..... imn
Day ..... id
Hour ..... ih
Minute ..... im
RAMS Grid Searched ..... ng
Number of sources ..... nsources
Total Particles ..... ntotparts
```

Followed by the source header lines (format *4i8*) - note that there can be only one species per source within the code:

```
Database Source I ..... (nsrc)
HYPACT internal Species ID ... pspecies(nsrc)
Database Species ID ..... pspecies(species(nsrc))
Source Particles ..... nsrcparts(nsrc)
```

Followed by one line for each particle (format *7e16.8,i5*):

```
x Location (m) ..... atp11
y Location (m) ..... atp12
z ASL (m) ..... atp13 + htopo
z AGL (m) ..... atp13
Mass of Particle ..... atp(1,8)
ppm Mass of Particle ..... atp(1,9)
Particle Age (s) ..... hytime - catp(1,1)
Hybrid Cluster ..... latp(1,2)
```

HYPACT Concentration and Header Files

VFILE Format Files (*a-C-*.g#* and *a-C-*.head*). Files output are limited to the grids selected by the [\\$OUTPUT namelist](#) parameter [irgrid](#).

Meteorological variables saved (if requested by [\\$OUTPUT namelist](#) parameter [imetout](#)):

- Topography
- Land Percentage
- U Wind Component
- V Wind Component
- W Wind Component
- Turbulent Kinetic Energy
- Theta
- Vertical Diffusivity
- Horizontal Diffusivity

Concentration variables saved for each species (if requested by [\\$OUTPUT namelist](#) parameters [ieulout](#) and [ilagout](#), and the [\\$SPECIES namelist](#) parameter [ioutspec](#)):

- Lagrangian
- Average Lagrangian
- Eulerian
- Average Lagrangian