



Part 9

Solid Modelling in Aspen Plus





Problem Definition

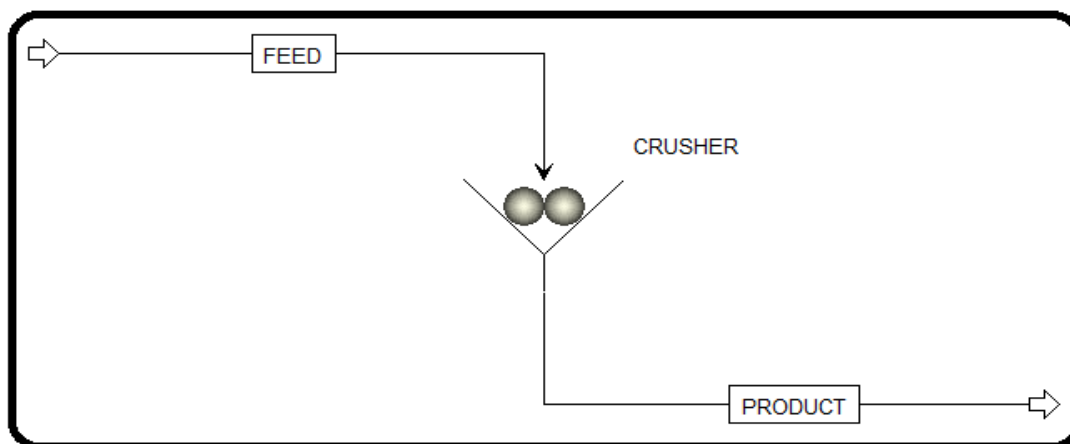
Solids handling is important in some chemical industries, such as specialty chemicals (e.g., fertilizers and silicones), extractive processes (e.g., oil shale, copper, and alumina), and biofuel processes (e.g., corn stover and sugarcane).

Unlike liquids, solids handling requires the knowledge of additional characteristics, such as average particle size and density, moisture content, solubility, color, shape, porosity, and particle size distribution, which is usually expressed in terms of the mean particle size and a factor describing the degree of scatter around the mean (e.g., standard deviation). Moreover, it is not possible to infer bulk from microscopic (single particles) properties of solids, in general. Aspen Plus has already many useful features that can deal with solids as it perfectly dealt with liquids and gases, as we have seen so far in the previous sessions. For more information about the different solids unit operations, refer to "APPENDIX".

Since there are many applications that involve more or less solid materials, we pick up some of them to demonstrate the important features of Aspen Plus in dealing with solid materials.

Example: THE CRUSHER

A feed stream of KCL with flowrate of 1250 kg/hr at 20 C and 1 bar with particle size distribution of median 5 mm and standard deviation of 1 mm with upper limit of 10 mm and lower limit of 0 mm passes a crusher which reduces the particle size of the feed stream.





How to simulate

1. Using Aspen Plus, start a new simulation by choosing the “Solids” category and selecting “Solids with Metric Units” template to create a steady-state flowsheet. Notice that Aspen Plus, does not assign a default property method. So, we will set it to “SOLIDS”.
2. In “Navigation” pane menu tree, go to “Components” | “Specifications” | “Selection” sheet. Click on “Find” button to search for “KCL”. Once you find “KCL”, add it to the list of components and change its “Type” from “Conventional” to “Solid”, as shown in Figure 14.2.

Component ID	Type	Component name	Alias	CAS number
KCL	Solid	POTASSIUM-CHLORIDE	KCL	7447-40-7
*				

The screenshot shows the Aspen Plus Properties window for the SOLIDS method. The window is divided into several sections:

- Global** (selected tab)
- Flowsheet Sections**
- Referenced**
- Comments**

Property methods & options

- Method filter: COMMON
- Base method: SOLIDS
- Henry components: (empty)

Petroleum calculation options

- Free-water method: STEAM-TA
- Water solubility: 3

Electrolyte calculation options

- Chemistry ID: (empty)
- ☒ Use true components

Method name: SOLIDS

Modify (checkbox checked)

- Vapor EOS: ESIG
- Data set: 1
- Liquid gamma: GMIDL
- Data set: 1
- Liquid molar enthalpy: HLMX108
- Liquid molar volume: VLMX25
- ☐ Heat of mixing
- ☐ Poynting correction
- ☒ Use liquid reference state enthalpy

NOTE #1: “KCL” being defined as solid type will be treated as an inert solid with known molecular structure and it will not be involved in aqueous phase equilibrium (i.e., no association/dissociation reactions). However, as we will see shortly that it requires knowledge of further attributes that have to deal with particle size distribution (PSD). Thus, the substream class type for the feed (or product) stream class will be Conventional Inert with PSD (“CIPSD”).



3. Click on “Next” button, run the simulator, and monitor serious warning and errors (if any) via the “Control Panel”. Once you successfully manage to complete the property analysis step, switch to “Simulation” environment.

4. From “Model Palette”, select “Solids” tab, click on “Crusher” icon, and add the crusher icon to the flowsheet area. In addition, add the proper input and output stream, as shown in Figure 14.1. Click on “Next” button and Aspen Plus will bring you to “FEED” stream input form.

Notice here that entering solid-containing feed stream properties are different from those of a conventional feed stream, which has no “Solid” type component. Figure 14.3 shows the typical input data to be entered for a feed stream; however, we will use the second tab, that is, “CI Solid” tab for a conventional inert solid as in our case.

Component	Value
KCL	1

Total 1

Notice that the half-filled red circle shown in Figure 14.3 indicates that input data pertaining to “FEED” stream is still incomplete, and the reason for this is simply because the stream class is of “MIXCIPSD” type with a substream of “CIPSD” type. Under such stream and substream definitions, we need to define or associate a PSD with the feed stream as well as define a simulation PSD for presentation of results.

5. Figure 14.4 (top) shows “Setup” | “Stream Class” | “Flowsheet” tab window where it shows that the stream class for the “Global” section (i.e., entire flowsheet) is of “MIXCIPSD” type. See “APPENDIX” for further information on stream classes. Stream classes ease the integration of solids and fluids in one simulation. The stream class can either be created or selected from a predefined type. Upon one’s need, a predefined class type can also be modified. Moreover, the “Stream Class” tab window (Figure 14.4 bottom) shows the selected substream class types for the given stream class of type “MIXCIPSD”. One can select the stream class from the drop-down list and see what substreams are already associated with.

Section	Stream class
GLOBAL	MIXCIPSD



Flowsheet Streams **Stream Class** Load Streams Comments

Stream class **MIXCIPSD**

Select substreams for stream class

CISOLID
NC
NCPSD

>
>>
<
<<

MIXED
CIPSD

< >
< >>
> <
>> <<

6. Defining the Particle Size Distribution (PSD)

Regarding the particle size distribution (PSD), we will define two PSD types one for the entire simulation and another for the feed stream. Notice here that both PSD types have to be within the same range; that is, the PSD mesh of the simulation within which we monitor the product stream solid attributes should not be too fine or too coarse for the feed stream PSD mesh. So, let us define the simulation PSD first in terms of size unit, lower limit, and upper limit. Figure 14.5 shows that we have selected the “PSD mesh type” to be “*Equidistant*” with parameters as shown in the figure. Do not forget to click on “Create PSD Mesh” button to confirm the creation of the selected PSD.

Mesh Comments

PSD mesh ID **PSD**

PSD Mesh

PSD mesh type **Equidistant**

No. of intervals **10**

Lower limit **0** Upper limit **10** Size units **mm**

Create PSD Mesh

Caution: If you reduce the number of intervals, data will be lost for the removed intervals wherever this PSD is used.

Particle size distribution mesh

	Int.	Lower	Upper
▶	1	0	1
▶	2	1	2
▶	3	2	3
▶	4	3	4
▶	5	4	5
▶	6	5	6
▶	7	6	7
▶	8	7	8
▶	9	8	9
▶	10	9	10

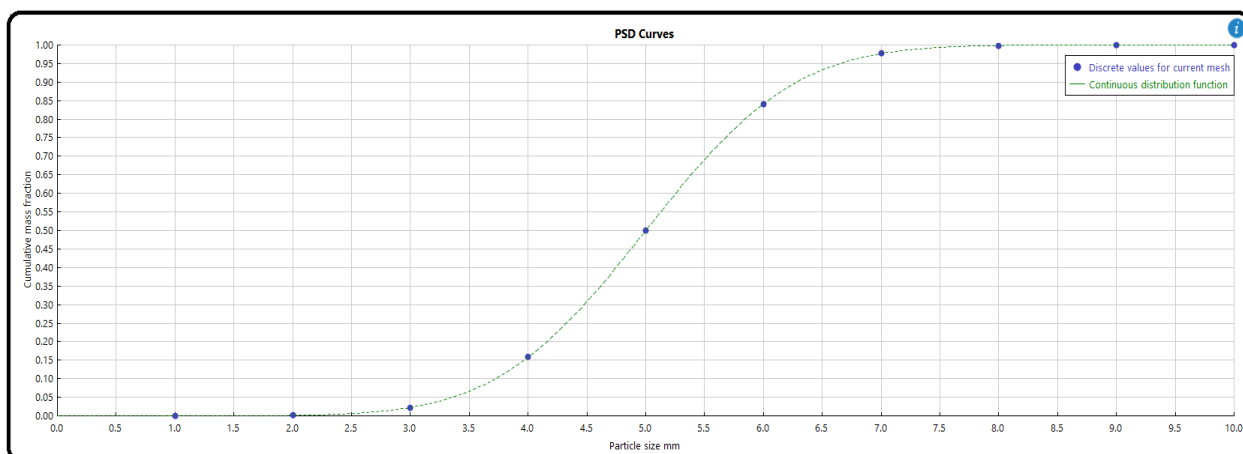


7. Figure 14.6 shows the “CI Solid” tab window for “FEED” stream where we need to define a PSD for the solid of the feed stream. Notice here that you can edit (or modify) the simulation PSD via clicking on “Edit PSD Mesh” button, which will invoke the form shown earlier in Figure 14.5. Moreover, we define a PSD of “FEED” stream, here, as “Normal” with a standard deviation of 1mm and median (D50) of 5 mm.

Interval	Lower limit	Upper limit	Weight fraction	Cumulative weight fraction
1	0	1	3.1686e-05	3.1686e-05
2	1	2	0.00131828	0.00134997
3	2	3	0.0214001	0.0227501
4	3	4	0.135905	0.158655
5	4	5	0.341345	0.5
6	5	6	0.341345	0.841345
7	6	7	0.135905	0.97725
8	7	8	0.0214001	0.99865
9	8	9	0.00131828	0.999969
10	9	10	3.13909e-05	1

Click on “Calculate” button to generate both the tabulated and graphic representation of PSD for “FEED” stream. Figure 14.7 shows the populated PSD values over the selected PSD mesh, which extends from 0 up to 10mm with equal subintervals. Notice that we have the frequency (i.e., weight fraction) is normally (i.e., symmetrically) distributed around its mean value that lies at about 5 mm. The range of simulation PSD acts like a window through which we have the chance to look at the effect of solids crushing on the PSD of the feed stream.

Figure 14.8 shows the plot of cumulative mass fraction as a function of particle size (mm).





8. Calculation of the Outlet PSD

Click on “Next” button and the “Blocks” | “CRUSHER” | “Input” | “Specifications” tab sheet will show up, as shown in Figure 14.9. The outlet PSD will be calculated based on one of three methods: The first method is to spell out the crusher type, breakage function parameters, impact/rotor velocity, and/or some of its sizing parameters.

The second method is based on known comminution (disintegration) power and some specific distribution function parameters. Finally, the third method is based on known outlet PSD, which can be either user-specified or common built-in distribution function, such as GGS, RRSB, Normal, and Log normal

The screenshot shows the 'Specifications' tab of a software interface for a CRUSHER block. The interface is organized into several sections:

- Outlet PSD calculation method:** Three radio buttons are present: 'Select equipment' (selected), 'Determine outlet PSD from comminution power and a distribution function', and 'Specify outlet PSD'.
- Operating parameters:** Four dropdown menus are shown: 'Crusher type' (Gyratory), 'Selection function' (US Bureau of Mines), 'Breakage function' (US Bureau of Mines), and 'Distribution function' (Rosin Rammler Sperling Bennet).
- Selection function parameters:** One dropdown menu is shown: 'Operating mode' (Primary).
- Breakage function parameters:** Three radio buttons are present: 'Maximum particle diameter' (selected), 'Power specification:', and 'Specific power:'. Below these are input fields for 'Maximum particle diameter' (2 mm), 'Power specification:' (kW), and 'Specific power:' (kWhr/ton). At the bottom, there is a field for 'Ratio of cut-off size to solids outlet diameter' (1.7).

9. Click on “Next” button, run the show, and watch out any serious warning or error in “Control Panel”. Figure 14.10 (left) shows the “Summary” tab window for “CRUSHER” block under “Results” sheet. Obviously, there is a size reduction between the feed and the outlet stream. The size reduction ratio is 5.656 evaluated at the median. The Sauter mean diameter (see “APPENDIX 14.E”) is also reduced from 4.76 to 0.725 mm. For example, Figure 14.10 (right) shows that 98.782% of total particles fall below 3mm diameter size or 99.99997% of them lie below the 4mm diameter dividing cut.



Summary	Balance	Utility Usage	✓ Status
Calculated power	1.17070888e-30	kW	▼
Particle diameter which is larger than 80% of inlet mass	0.00587888	meter	▼
Particle diameter which is larger than 80% of outlet mass	0.00183233	meter	▼
Particle diameter which is larger than 50% of inlet mass	0.005	meter	▼
Particle diameter which is larger than 50% of outlet mass	0.00088407	meter	▼
Size reduction ratio of D80	3.20841		
Size reduction ratio of D50	5.65566		
Sauter mean diameter of inlet particles	0.00475806	meter	▼
Sauter mean diameter of outlet particles	0.000725362	meter	▼



10. Repeat the same task of reducing the particle size of “FEED” stream but this time using the second method, which is “*Determine outlet PSD from comminution power and a distribution function*” (see Figure 14.9). Under “Specifications” tab, use “*Rosin Rammler Sperling Bennet*” distribution function; power specification of 1kW; and select “*D50*” parameter with “*1mm*” from “Distribution function parameters” option. Under “Grindability” tab, use “*Bond’s law*” as the “Comminution law” and for “CIPSD” substream use “*3 kWhr/ton*” under “Bond work index” column. See Aspen Plus built-in help: “Bond’s Law” for further information on Bond work index. Bond considered that the work necessary for reduction is inversely proportional to the square root of the size produced. This applies for particles between 0.05 and 50 mm.



Summary	Balance	Utility Usage	Status
Calculated power	1	kW	
Particle diameter which is larger than 80% of inlet mass	0.00587888	meter	
Particle diameter which is larger than 80% of outlet mass	0.0016	meter	
Particle diameter which is larger than 50% of inlet mass	0.005	meter	
Particle diameter which is larger than 50% of outlet mass	0.000999989	meter	
Size reduction ratio of D80	3.67431		
Size reduction ratio of D50	5.00006		
Sauter mean diameter of inlet particles	0.00475806	meter	
Sauter mean diameter of outlet particles	0.000749996	meter	

Material	Work	Vol.% Curves	Wt. % Curves	Petroleum	Polymers	Solids
			Units	FEED	PRODUCT	
+	Mass Fractions					
	Volume Flow		cum/hr	0.628824	0.628824	
-	PSD					
	0 - 1 mm			3.1686e-05	0.500006	
	- 2 mm			0.00131828	0.499994	
	- 3 mm			0.0214001	0	
	- 4 mm			0.135905	0	
	- 5 mm			0.341345	0	
	- 6 mm			0.341345	0	
	- 7 mm			0.135905	0	
	- 8 mm			0.0214001	0	
	- 9 mm			0.00131828	0	
	- 10 mm			3.13989e-05	0	

11. Repeat the same task of reducing the particle size of "FEED" stream but this time using the third method, that is, "Specify outlet PSD" (see Figure 14.9). Under "Specifications" tab, select "Substream ID" as CIPSD; select "Use distribution function"; the "Bypass fraction" is set to 0; create a new distribution function ID and name it 1; select the "Distribution function" as Normal; select D50 parameter with 1mm and standard deviation of 1 mm.

Under "Grindability" tab, you may keep it blank (i.e., no entries at all); however, if you use Bond's law as the comminution law and for "CIPSD" substream use "3 kWhr/ton" under "Bond work index"



column, then you will be able to calculate the power needed to achieve the mission, that is, size reduction of feed stream particles down to the specified outlet PSD.

Report the outlet stream solid attributes in terms of particle size cut, reduction, mean, and distribution (i.e., PSD). Also, report the calculated power (kW) needed to carry out the crushing mission. I hope toward the end of this exercise that you managed to crush the information needed to deal with crushers.

Summary	Balance	Utility Usage	✓ Status
Calculated power	0.555457537	kW	▼
Particle diameter which is larger than 80% of inlet mass	0.00587888	meter	▼
Particle diameter which is larger than 80% of outlet mass	0.00187888	meter	▼
Particle diameter which is larger than 50% of inlet mass	0.005	meter	▼
Particle diameter which is larger than 50% of outlet mass	0.001	meter	▼
Size reduction ratio of D80	3.12893		
Size reduction ratio of D50	5		
Sauter mean diameter of inlet particles	0.00475806	meter	▼
Sauter mean diameter of outlet particles	0.000776194	meter	▼

[illegible]



Part 2

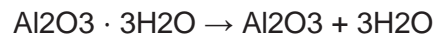
Problem Definition

THE FLUIDIZED BED FOR ALUMINA DEHYDRATION

In this problem, we will handle the fluidized bed with a chemical reaction. Fluidized beds are used in a variety of industrial processes, which include drying, cooling, heating, and as reactors. The feed stream is made of aluminum trihydroxide ($\text{Al}(\text{OH})_3$), which will be dispensed in the form of solid particles to the entrance of the fluidized bed where it will be pneumatically lifted or fluidized by means of a hot air stream. This will cause the phenomenon known as fluidization visualizing solid particles behaving as fluid-like molecules (i.e., they gain the translational, vigorous vibrational, and rotational kinetic energy as well). This being the case, we will enhance the simultaneous mass and heat transfer rates from and to the solid particles. The dehydration reaction



can be put in another form:



So, the reaction is basically a solid transformation from the hydrated into the anhydrous unit crystal. Thus, heating by convection (by the air) and conduction (through the solid particle) will liberate the combined water molecules from the crystal. Keep in mind that this process is a sort of “roasting” process rather than drying. For a typical wet-hydrated alumina or inorganic solids, in general, water molecules are trapped within the solid matrix itself (i.e., within the macropores of a solid pellet, compared with the tiny small size of the unit crystal).



How to Simulate

1. Using Aspen Plus, start a new simulation by choosing the “Solids” category and selecting “Solids with Metric Units” template to create a steady-state flowsheet. We will set the default property method to “SOLIDS”.
2. In “Navigation” pane menu tree, go to “Components” | “Specifications” | “Selection” tab sheet. Add the following components and change the “Type” of both “ $AL(OH)_3$ ” and “ AL_2O_3 ” from “Conventional” to “Solid”, as shown in Figure 14.11.

The screenshot shows the 'Selection' tab in the Aspen Plus interface. The 'Component ID' column lists AL(OH)3, AL2O3, AIR, and WATER. The 'Type' column shows Solid for AL(OH)3 and AL2O3, and Conventional for AIR and WATER. The 'Component name' column lists ALUMINIUM-HYDROXIDE, ALUMINIUM-OXIDE-ALPHA-C..., AIR, and WATER. The 'Alias' column lists AL(OH)3, AL2O3, AIR, and H2O. The 'CAS number' column lists 21645-51-2, 1344-28-1, 132259-10-0, and 7732-18-5.

Component ID	Type	Component name	Alias	CAS number
AL(OH)3	Solid	ALUMINIUM-HYDROXIDE	AL(OH)3	21645-51-2
AL2O3	Solid	ALUMINIUM-OXIDE-ALPHA-C...	AL2O3	1344-28-1
AIR	Conventional	AIR	AIR	132259-10-0
WATER	Conventional	WATER	H2O	7732-18-5

The screenshot shows the 'Global' tab in the Aspen Plus interface. The 'Method filter' is set to COMMON, the 'Base method' is set to SOLIDS, and the 'Henry components' are empty. The 'Petroleum calculation options' section shows 'Free-water method' set to STEAM-TA and 'Water solubility' set to 3. The 'Electrolyte calculation options' section shows 'Chemistry ID' as empty and 'Use true components' checked. The 'Method name' is set to SOLIDS. The 'Modify' section shows 'Vapor EOS' set to ESIG, 'Data set' set to 1, 'Liquid gamma' set to GMIDL, 'Data set' set to 1, 'Liquid molar enthalpy' set to HLMX108, 'Liquid molar volume' set to VLMX25, 'Heat of mixing' unchecked, 'Poynting correction' unchecked, and 'Use liquid reference state enthalpy' checked.

Property methods & options

Method filter: COMMON

Base method: SOLIDS

Henry components:

Petroleum calculation options

Free-water method: STEAM-TA

Water solubility: 3

Electrolyte calculation options

Chemistry ID:

☒ Use true components

Method name: SOLIDS

Methods Assistant...

☐ Modify

Vapor EOS: ESIG

Data set: 1

Liquid gamma: GMIDL

Data set: 1

Liquid molar enthalpy: HLMX108

Liquid molar volume: VLMX25

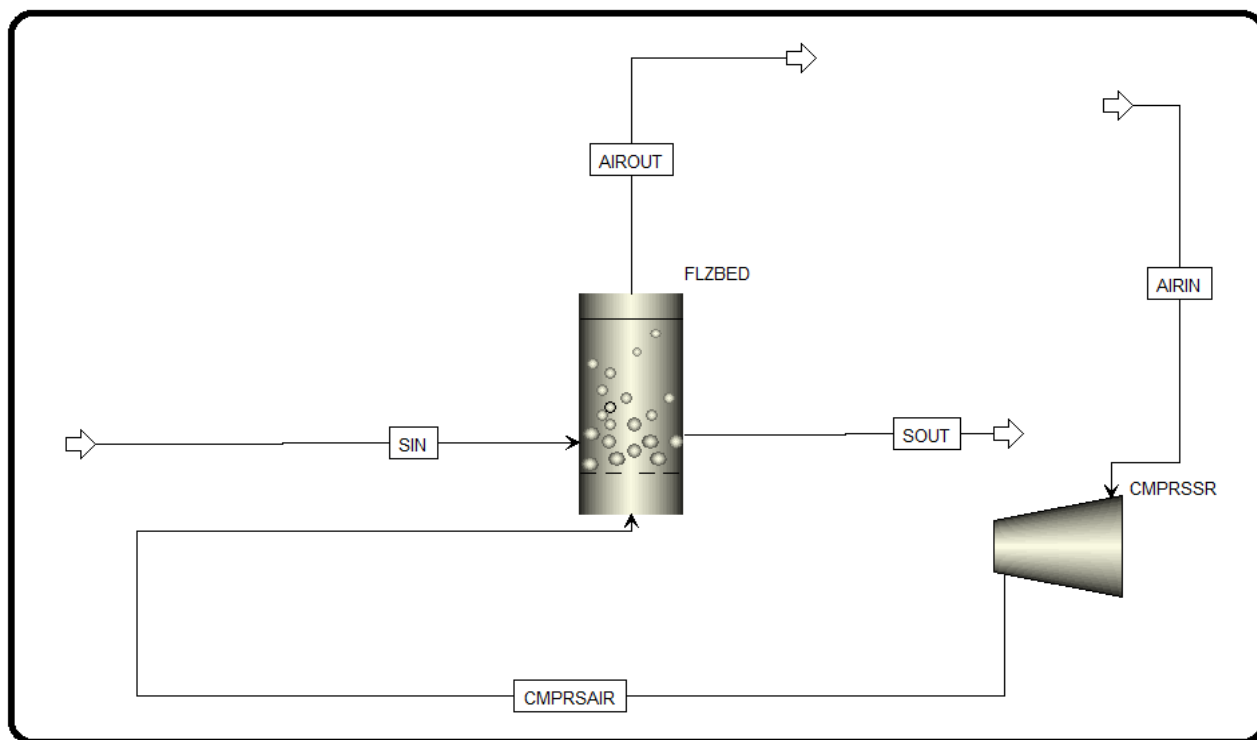
☐ Heat of mixing

☐ Poynting correction

☒ Use liquid reference state enthalpy



3. Click on “Next” button, run the simulator, and monitor warnings and errors (if any) via the “Control Panel”. Once you successfully manage to complete the property analysis step, switch to “Simulation” environment.
4. From “Model Palette”, select “Solids” tab, click on “Fluidbed” icon, and add the fluidized bed icon to the flowsheet area. Add one compressor unit from “Pressure Changers” tab. In addition, add the proper input and output streams, as shown in Figure 14.12. Notice that “SIN” and “SOUT” stream stand for the inlet and outlet solid streams, respectively.



5. Air stream (“**AIRIN**”) will be introduced to the compressor at a rate of 4500 kg/h at room temperature and pressure and the discharge pressure of the compressor will be set at 12 bar. Of course, the temperature of air will increase as a result of adiabatic compression. On the other hand, the solid feed stream (“**SIN**”) will enter at a rate of 300 kg/h at room temperature and a pressure of 12 bar as pure $\text{Al}(\text{OH})_3$. Moreover, Figure 14.13 (*left*) shows the simulation PSD mesh and that of “**SIN**” stream (*right*).



Mixed

CI Solid

NC Solid

Flash Options

EO Options

Costing

Comments

Specifications

Flash Type

Temperature

Pressure

State variables

Temperature

25

C

Pressure

1

bar

Vapor fraction

Total flow basis

Mass

Total flow rate

4500

kg/hr

Solvent

Reference Temperature

Volume flow reference temperature

C

Component concentration reference temperature

C

Composition

Mass-Frac

Component	Value
AL(OH)3	
AL2O3	
AIR	1
WATER	

Total1



Mixed

CI Solid

NC Solid

Flash Options

EO Options

Costing

Comments

Specifications

State variables

Substream name

CIPSD

Temperature

25

C

Pressure

12

bar

Total flow basis

Mass

Total flow rate

300

kg/hr

Composition

Mass-Flow

kg/hr

Component	Value
AL(OH)3	1
AL2O3	
Total	1

Component Attribute

Particle Size Distribution

PSD mesh ID

PSD

Units

mm

Edit PSD Mesh

Populate PSD using

User-specified values

☒ A distribution function

Distribution function

Type

GGS

Dispersion parameter

1.5

Maximum diameter

10

mm

Calculate



☒ Mesh
 ☐ Comments

PSD mesh ID

PSD Mesh

PSD mesh type

No. of intervals

Lower limit
 Upper limit
 Size units

Caution: If you reduce the number of intervals, data will be lost for the removed intervals wherever this PSD is used.

Particle size distribution mesh

	Int.	Lower	Upper
▶	1	0	1
▶	2	1	2
▶	3	2	3
▶	4	3	4
▶	5	4	5
▶	6	5	6
▶	7	6	7
▶	8	7	8
▶	9	8	9
▶	10	9	10

☒ Specifications
 ☐ Calculation Options
 ☐ Power Loss
 ☐ Convergence
 ☐ Integration Parameters
 ☐ Utility
 ☐ Comments

Model and type

Model ☒ Compressor ☐ Turbine

Type

Outlet specification

☒ Discharge pressure

☐ Pressure increase

☐ Pressure ratio

☐ Power required

☐ Use performance curves to determine discharge conditions

Efficiencies

Isentropic
 Polytropic
 Mechanical

6. For the fluidized bed block, the following design specifications are entered, as shown in Figure 14.14. First, under “Specifications” tab, the bed mass is set to *120 kg*; “*Geldart B*” classification; and “*Ergun*” model for finding the minimum fluidization velocity. Second, under “Geometry” tab, the bed height is set to *4m*; the solid discharge location *0.95* (i.e., $0.95 \times 4 = 3.8\text{m}$, measured from the bottom of the bed); and the cross-sectional area with a constant diameter equal to *0.5m*. Third, under “Gas Distributor” tab, its type is set to “*Perforated plate*” with number of orifices equal to *40*, diameter equal to *10mm* each, and the default orifice discharge coefficient of *0.8*. Fourth, under “Convergence” tab, for mass balance convergence, the solver is set to “*Newton*” instead of



“Broyden”, as the latter did not properly converge for a reasonable solution for the given fluidization case. Finally, under “Reactions” tab we will associate the reaction “R-1” set to the fluidized bed. The reaction “R-1” set will be created shortly.

Specifications	Operation	Geometry	Gas Distributor	Heat Exchanger	Reactions	PSD	Convergence	Comments
Bed inventory								
<input checked="" type="radio"/> Specify bed mass		120		kg				
<input type="radio"/> Specify bed pressure drop				bar				
Voidage at minimum fluidization		0.5						
Geldart classification		Geldart B						
Minimum fluidization velocity								
<input type="radio"/> Specify velocity				m/sec				
<input checked="" type="radio"/> Calculate from correlation		Ergun						
Transport disengagement height								
TDH model		George and Grace						
Maximum dCv/dh		1e-05						
Elutriation								
Model		Tasirin & Geldart						
Decay constant		3						
TG parameter A1		23.7		$k_{i,\infty} = A \cdot \rho_G \cdot u^B \cdot \exp\left(C - \frac{u_{t,i}}{u}\right)$				
TG parameter A2		14.5						

Specifications	Operation	Geometry	Gas Distributor	Heat Exchanger	Reactions	PSD	Convergence
Dimensions							
Height		4		meter			
Solids discharge location		0.95					
Cross-section		Circular					
<input checked="" type="radio"/> Constant diameter		0.5		meter			
<input type="radio"/> Height-dependent diameter							
Width				meter			
Depth				meter			
Height-dependent diameter							
Location		Diameter					
		meter					
Secondary gas inlet streams							
Name		Location					



Specifications Operation Geometry Gas Distributor Heat Exchanger Reactions PSD Convergence

Type *Perforated plate*

Number of orifices 40

Orifice diameter 10 mm

☒ Specify orifice discharge coefficient 0.8

☐ Specify distributor pressure drop

Number of bubble caps

Number of cap orifices

Cap orifice diameter meter

☐ Specify cap orifice discharge coefficient 0.8

☐ Specify distributor pressure drop bar

Specifications Operation Geometry Gas Distributor Heat Exchanger Reactions PSD Convergence

Select reaction sets to be included in the model

Available reaction sets	Selected reaction sets
	R-1

> >> < << New

Specifications Operation Geometry Gas Distributor Heat Exchanger Reactions PSD Convergence

PSD calculation option

☒ Keep PSD

☐ Calculate PSD from particle growth model

☐ User-specified PSD



Specifications Operation Geometry Gas Distributor Heat Exchanger Reactions PSD Convergence

Convergence parameters

Tolerance 0.0001

Maximum number of iteration steps 100

Number of cells for the bottom zone 100

Number of cells for the dilute zone 100

Bed mass threshold 1e-10

Pressure drop threshold 1e-10

Holdup convergence method RootN1

Pressure convergence method RootN1

Flash options

Tolerance 0.0001

Maximum number of iterations 30

Temperature estimate C

Reaction convergence

Mass balance convergence

Solver Newton Newton Parameters

The reaction “R-1” set will be created under “Reactions” folder. Figure 14.15 shows the stoichiometry and kinetic data associated with alumina transformation from a hydrous to anhydrous state.

Stoichiometry Kinetic Equilibrium Activity Comments				
New Edit Copy Paste				
Rxn No.	Reaction type	Stoichiometry		Delete
1	Kinetic	2 AL(OH)3(CIPSD) --> AL2O3(CIPSD) + 3 WATER(MIXED)		X



Edit Reaction

Reaction No. ☒ 1

Reaction type *Kinetic*

Reactants

	Component	Coefficient	Exponent
▶	AL(OH)3 (CIPSD)	-2	1
▶			

Products

	Component	Coefficient	Exponent
▶	AL2O3 (CIPSD)	1	
▶	WATER	3	
▶			

☒ Stoichiometry ☒ **Kinetic** Equilibrium Activity Comments

1) 2 AL(OH)3(CIPSD) --> AL2O3(CIPSD) + 3 WATER(MIXED)

Reacting phase **Vapor** Rate basis *Reac (vol)*

Power Law kinetic expression

If To is specified Kinetic factor $=k(T/T_o)^n e^{-(E/R)[1/T-1/T_o]}$

If To is not specified Kinetic factor $=kT^n e^{-E/RT}$

k

n

E

To

[Ci] basis *Molarity*



Results

Figure 14.16 (*left*) shows the results summary of the calculated geometrical and operational variables of the fluidized bed and the moisture profile of air stream where dry air enters at the bottom and becomes moist as it moves upward (*right*).

Summary	Balance	Profiles	Gas compositions	Status
Height of bottom zone	0.108436	meter		
Height of freeboard	3.89156	meter		
TDH calculated from correlation	2.19969	meter		
TDH based on solids volume fraction profile	3.85265	meter		
Solids holdup	119.998	kg		
Number of particles in bed	1.60882e+07			
Surface area	48.3044	sqm		
Distributor pressure drop	0.211705	bar		
Bottom zone pressure drop	0.0154238	bar		
Freeboard pressure drop	0.0469366	bar		
Fluidized bed pressure drop	0.0623604	bar		
Overall pressure drop	0.274065	bar		
Heat duty	0	Gcal/hr		
Minimum fluidization velocity	0.704601	m/sec		
Calculated temperature	353.468	C		
Moles generated	3.84505	kmol/hr		



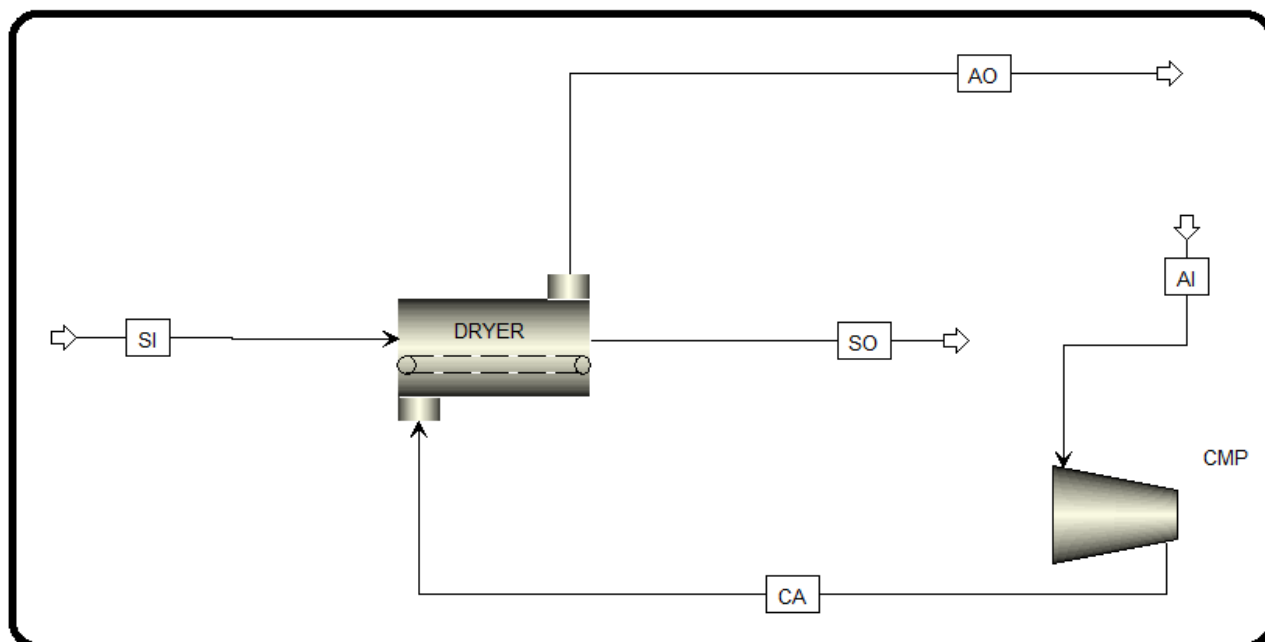
Material	Vol.% Curves	Wt. % Curves	Petroleum	Polymers	Solids					
			Units	CMPSAIR	SIN	AIROUT	SOUT			
▶	— Total Stream									
▶	Temperature	C	442.065	25	353.468	353.468				
▶	Pressure	bar	12	12	11.7259	11.7261				
▶	Molar Vapor Fraction		1	0	0.999999	0				
▶	Molar Liquid Fraction		0	0	0	0				
▶	Molar Solid Fraction		0	1	5.92826e-07	1				
▶	Mass Vapor Fraction		1	0	0.999998	0				
▶	Mass Liquid Fraction		0	0	0	0				
▶	Mass Solid Fraction		0	1	2.11621e-06	1				
▶	Molar Enthalpy	kcal/mol	2.97075	-304.767	0.267563	-391.763				
▶	Mass Enthalpy	kcal/kg	102.614	-3907.1	9.36854	-3842.71				
▶	Molar Entropy	cal/mol-K	1.292	-96.9643	0.325636	-56.6278				
▶	Mass Entropy	cal/gm-K	0.0446271	-1.24308	0.0114019	-0.555448				
▶	Molar Density	mol/cc	0.000201798	0.031024	0.00022507	0.0390551				



Part 3

Problem Definition: KCL DRYING

Consider the process of KCl drying using hot air, where moisture will migrate from the wet solid to the dry air and the heat needed will be convectively supplied by the hot air.





How to Simulate

1. Using Aspen Plus, start a new simulation by choosing the “Solids” category and selecting “Solids with Metric Units” template to create a steady-state flow sheet. Notice that Aspen Plus does not assign a default property method. So, we will set it to “SOLIDS”.
2. In “Navigation” pane menu tree, go to “Components” | “Specifications” | “Selection” sheet. Click on “Find” button to search for “KCL”. Once you find “KCL”, add it to the list of components and change its “Type” from “Conventional” to “Solid”. Add water and air as well, as shown in Figure 14.18.

Component ID	Type	Component name	Alias	CAS number
KCL	Solid	POTASSIUM-CHLORIDE	KCL	7447-40-7
AIR	Conventional	AIR	AIR	132259-10-0
H2O	Conventional	WATER	H2O	7732-18-5
*				

Buttons: Find, Elec Wizard, SFE Assistant, User Defined, Reorder, Review

Property methods & options

Method filter: COMMON

Base method: SOLIDS

Henry components: [Empty]

Petroleum calculation options

Free-water method: STEAM-TA

Water solubility: 3

Electrolyte calculation options

Chemistry ID: [Empty]

☒ Use true components

Method name: SOLIDS

Methods Assistant...

☐ Modify

Vapor EOS: ESIG

Data set: 1

Liquid gamma: GMIDL

Data set: 1

Liquid molar enthalpy: HLMX108

Liquid molar volume: VLMX25

☐ Heat of mixing

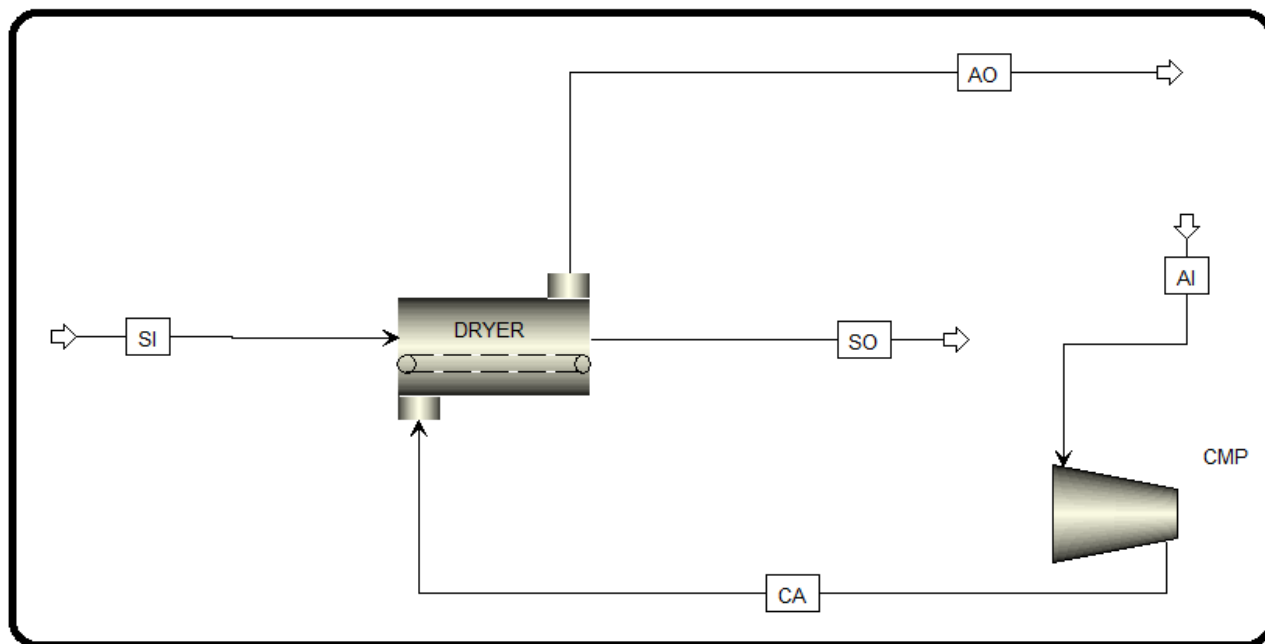
☐ Poynting correction

☒ Use liquid reference state enthalpy

3. Click on “Next” button, run the simulator, and monitor warnings and errors (if any) via the “Control Panel”. Once you successfully manage to complete the property analysis step, switch to “Simulation” environment. From “Model Palette”, select “Solids” tab, click on “Dryer” icon, and add



the dryer icon to the flowsheet area. Add one compressor unit from “Pressure Changers” tab. In addition, add the proper input and output streams, as shown in Figure 14.19.



Under “**Solids**” folder, select water as the moisture component, as shown in Figure 14.20.

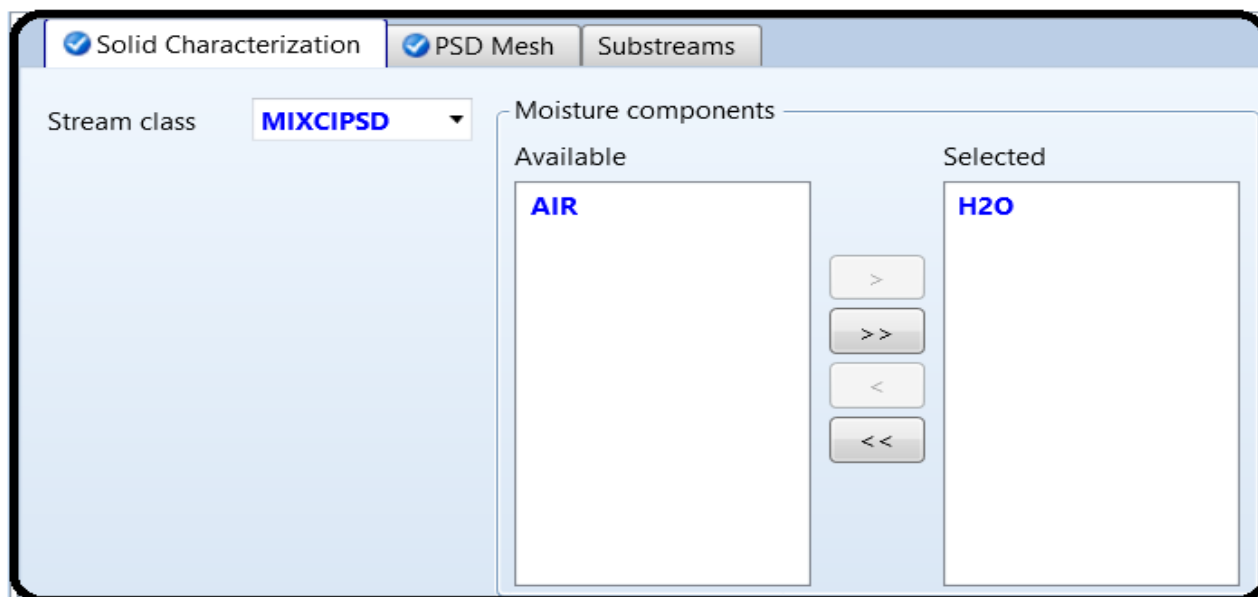


Figure 14.21 shows the creation of a simulation PSD mesh for monitoring the effect of drying on the particle size distribution of the substance being dried.



Mesh | Comments

PSD mesh ID: PSD

PSD Mesh

PSD mesh type: **Equidistant**

No. of intervals: **10**

Lower limit: **1** Upper limit: **5** Size units: **cm**

Create PSD Mesh

Caution: If you reduce the number of intervals, data will be lost for the removed intervals wherever this PSD is used.

Particle size distribution mesh

	Int.	Lower	Upper
	1	1	1.4
	2	1.4	1.8
	3	1.8	2.2
	4	2.2	2.6
	5	2.6	3
	6	3	3.4
	7	3.4	3.8
	8	3.8	4.2
	9	4.2	4.6
	10	4.6	5

Figure 14.22 shows the solid inlet stream properties in terms of P , T , compositional flowrate, and PSD

Mixed | **CI Solid** | NC Solid | Flash Options | EO Options | Costing | Comments

Specifications

State variables

Substream name: **CIPSD**

Temperature: **25 C**

Pressure: **1 bar**

Total flow basis: **Mass**

Total flow rate: **100 kg/hr**

Composition

Mass-Frac

Component	Value
KCL	0.8
H2O	0.2
Total	1

Component Attribute

Particle Size Distribution

PSD mesh ID: **PSD** Units: **cm**

Edit PSD Mesh

Populate PSD using

☐ User-specified values

☒ A distribution function

Distribution function

Type: **Normal**

Standard deviation: **1 cm**

D50: **2.4 cm**

Calculate



The inlet air stream enters the compressor at a flow rate of 3000 kg/h at room temperature and pressure and will leave it at 6 bar. The dryer specifications are shown in Figure 14.23.

Mixed

CI Solid

NC Solid

Flash Options

EO Options

Costing

Comments

Specifications

Flash Type

Temperature

Pressure

State variables

Temperature

25

C

Pressure

1

bar

Vapor fraction

Total flow basis

Mass

Total flow rate

3000

kg/hr

Solvent

Reference Temperature

Volume flow reference temperature

C

Component concentration reference temperature

C

Composition

Mass-Frac

Component	Value
KCL	
AIR	1
H2O	

Total

1

5. Click Next to go to Compressor Block.



Specifications Calculation Options Power Loss Convergence Integration Parameters

Model and type

Model ☒ Compressor ☐ Turbine

Type **Isentropic**

Outlet specification

☒ Discharge pressure **bar**

☐ Pressure increase bar

☐ Pressure ratio

☐ Power required kW

☐ Use performance curves to determine discharge conditions

Efficiencies

Isentropic Polytropic Mechanical

Notice that the “**Shortcut**” is simpler than the “**Convective dryer**” method in terms of input requirements, nevertheless, less rigorous. The drying curve characteristics are shown in Figure 14.24. Notice that the **critical solids moisture content** is the moisture content at which further evaporation is mass-transfer limited and the second drying phase begins, and the **equilibrium solids moisture content** is the ultimate moisture content at which point no further drying is possible.

Specifications PSD Entrainment ☒ Mass/Heat Transfer Atomization ☒ Drying Curve Convergence Utility Comments

Operation mode **Continuous**

Dryer type: **Convective dryer**

Geometry

Gas flow direction: **Cross-flow**

Solids flow: **Plug flow**

Input specifications: **Length** **Solids residence time**

Volume: cum

Length: **meter**

Cross sectional area: sqm

Solids residence time: **min**

Solids velocity: m/sec

Solids holdup: kg

Fill grade:

Bed porosity:



Results

Summary	Balance	Evaporation	Profiles	Evaporation Rate	Utility Usage	Status
Exhaust gas temperature	276.304908	C				
Calculated duty	1.4428e-08	Gcal/hr				
Solids residence time	0.025	hr				
Solids velocity	0.0444444	m/sec				
Critical moisture content (dry)	0.0526316					
Equilibrium moisture content (dry)	0.0309278					
Initial solids moisture content (dry)	0.25					
Outlet solids moisture content (dry)	0.0309278					
Overall evaporation rate	17.5258	kg/hr				
Initial inlet vapor moisture content (dry)	0					
Outlet vapor moisture content (dry)	0.00584192					
Vapor temp at adiabatic saturation	53.9143	C				
Vapor moisture content at adiabatic saturation	0.109343					
Calculated Lewis number	0.805167					
Calculated heat transfer coefficient	1502.07	kcal/hr-sqm-K				

Material	Heat	Load	Vol.% Curves	Wt. % Curves	Petroleum	Polymers	Solids				
				Units	CA	SI	AO	SO			
	▶	Mass Density	kg/cum		3.65269	1656.38	0.631502	1835.84			
	▶	Enthalpy Flow	Gcal/hr		0.19935	-0.187651	0.128439	-0.116741			
	▶	Average MW			28.9509	45.8031	28.8492	68.1362			
	▶	+ Mole Flows	kmol/hr		103.624	2.18326	104.597	1.21043			
	▶	+ Mole Fractions									
	▶	- Mass Flows	kg/hr		3000	100	3017.53	82.4742			
	▶	KCL	kg/hr		0	80	0	80			
	▶	AIR	kg/hr		3000	0	3000	0			
	▶	H2O	kg/hr		0	20	17.5258	2.47423			
	▶	+ Mass Fractions									
	▶	Volume Flow	cum/hr		821.313	0.0603725	4778.33	0.0449245			



Example: KCL CRYSTALLIZATION

Consider the process of KCl crystallization from KCl solution using heat to evaporate the water and result in a supersaturated solution that will be the cause for solid precipitation at the bottom of the mother liquor.

1. Using Aspen Plus, start a new simulation by choosing the “Solids” category and selecting “Solids with Metric Units” template to create a steady-state flowsheet. Notice that Aspen Plus does not assign a default property method. So, we will set it to “SOLIDS”.
2. In “Navigation” pane menu tree, go to “Components” | “Specifications” | “Selection” sheet. Add KCl twice once as “Solid” (i.e., “KCL(S)”) and another as “Conventional” (i.e., “KCL”) to the list of components and add water as well, as shown in Figure 14.25.

The screenshot shows the 'Selection' sheet in Aspen Plus. The 'Component ID' column lists KCL(S), H2O, and KCL. The 'Type' column lists Solid, Conventional, and Conventional. The 'Component name' column lists POTASSIUM-CHLORIDE, WATER, and POTASSIUM-CHLORIDE. The 'Alias' column lists KCL, H2O, and KCL. The 'CAS number' column lists 7447-40-7, 7732-18-5, and 7447-40-7. The 'Find' button is highlighted.

Component ID	Type	Component name	Alias	CAS number
KCL(S)	Solid	POTASSIUM-CHLORIDE	KCL	7447-40-7
H2O	Conventional	WATER	H2O	7732-18-5
KCL	Conventional	POTASSIUM-CHLORIDE	KCL	7447-40-7

The screenshot shows the 'Global' sheet in Aspen Plus. The 'Method filter' is set to COMMON. The 'Base method' is set to SOLIDS. The 'Henry components' are empty. The 'Petroleum calculation options' section shows 'Free-water method' set to STEAM-TA and 'Water solubility' set to 3. The 'Electrolyte calculation options' section shows 'Chemistry ID' empty and 'Use true components' checked. The 'Method name' is set to SOLIDS. The 'Modify' section shows 'Vapor EOS' set to ESIG, 'Data set' set to 1, 'Liquid gamma' set to GMIDL, 'Data set' set to 1, 'Liquid molar enthalpy' set to HLMX108, 'Liquid molar volume' set to VLMX25, 'Heat of mixing' unchecked, 'Poynting correction' unchecked, and 'Use liquid reference state enthalpy' checked.

Property methods & options

Method filter: COMMON

Base method: SOLIDS

Henry components:

Petroleum calculation options

Free-water method: STEAM-TA

Water solubility: 3

Electrolyte calculation options

Chemistry ID:

☒ Use true components

Method name: SOLIDS

☐ Modify

Vapor EOS: ESIG

Data set: 1

Liquid gamma: GMIDL

Data set: 1

Liquid molar enthalpy: HLMX108

Liquid molar volume: VLMX25

☐ Heat of mixing

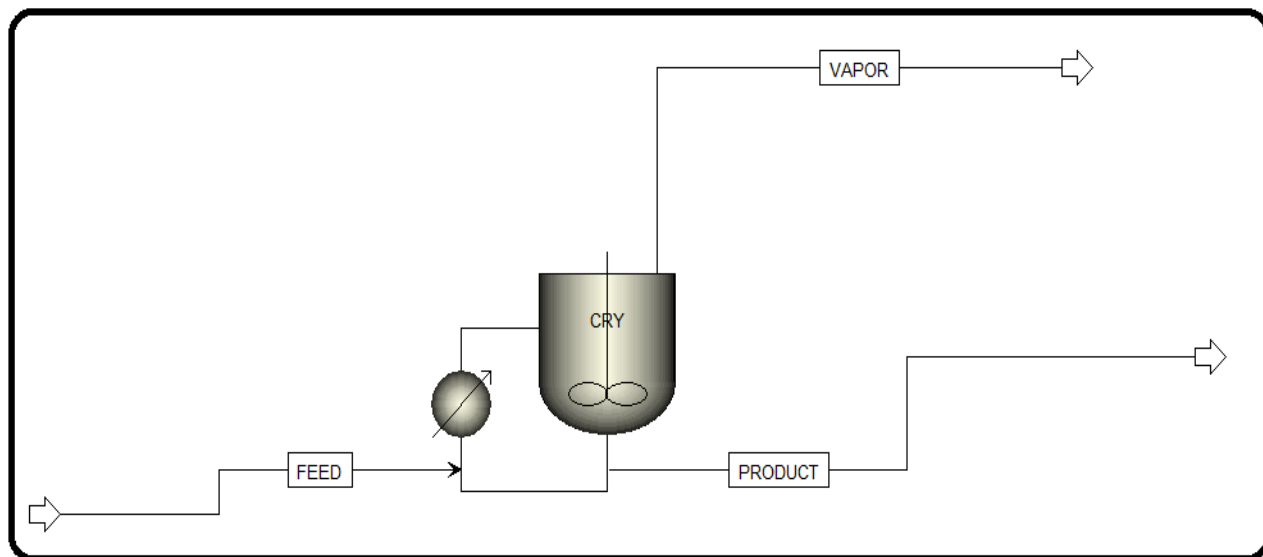
☐ Poynting correction

☒ Use liquid reference state enthalpy

3. Click on “Next” button, run the simulator, and monitor warning and errors (if any) via the “Control Panel”. Once you successfully manage to complete the property analysis step, switch to “Simulation” environment.



4. From “Model Palette”, select “Solids” tab, click on “Crystallizer” icon, and add the crystallizer icon to the flowsheet area. In addition, add the proper input and output streams, as shown in Figure 14.26.



5. Figure 14.29 shows the feed stream properties in terms of P , T , and compositional flowrate. There will be no need for PSD specification because the feed stream is a saturated aqueous KCl solution entering at the maximum solubility, which is 35 g KCl/100 g water (amounts to KCl mass fraction of 0.26).

Component	Value
KCL(S)	
H2O	0.74
KCL	0.26

Total 1



Under “**Solids**” folder, select water as the moisture component, as shown in Figure 14.27.

Stream class: **MIXCIPSD**

Moisture components

Available	Selected
KCL	H2O

Navigation buttons: >, >>, <, <<

PSD mesh ID: PSD

PSD Mesh

PSD mesh type: **Equidistant**

No. of intervals: **10**

Lower limit: **0** Upper limit: **5** Size units: **cm**

Create PSD Mesh

Caution: If you reduce the number of intervals, data will be lost for the removed intervals wherever this PSD is used.

	Int.	Lower	Upper
>	1	0	0.5
>	2	0.5	1
>	3	1	1.5
>	4	1.5	2
>	5	2	2.5
>	6	2.5	3
>	7	3	3.5
>	8	3.5	4
>	9	4	4.5
>	10	4.5	5

Figure 14.30 shows the “**Specifications**” tab window for the crystallizer. A heat duty of 300kW will be supplied to evaporate water (the solvent) and concentrate KCl solution. The “**Operating mode**” is set to “*Crystallizing*”. Under this mode, the amount of crystals should increase in the flow and a warning will be issued if the crystal product flow rate is zero or smaller than that of the inlet. The opposite is true for the “*Dissolving or melting*” mode. No warning will be issued if the



mode is set to “Either”. Figure 14.31 shows the “Crystallization” tab window where we define the sort of speaking “physical” reaction: $\text{KCl(aq)} \rightarrow \text{KCl(S)}$.

Operating conditions

Pressure: 1 bar

Heat duty: 300 kW

Saturation calculation method

- ☒ Solubility data
- ☐ Solubility function
- ☐ Chemistry
- ☐ User subroutine

Salt specifications

Salt component ID: [dropdown]

- ☒ Move crystallizing salt from Cl Solids substream
- ☒ Move crystallized salt to Cl Solids substream

Valid phases: Vapor-Liquid

Operating mode: Crystallizing

Reactants

Component	Coefficient
KCL	-1

Crystal product

KCL(S) (CIPSD)

Coefficient: 1

Stoichiometry

KCL --> KCL(S) (CIPSD)



Figure 14.32 shows the “Solubility” tab window where we enter the solubility data for KCl in water at room temperature. In the literature, it is reported as 35 g KCl per 100 g H₂O.

Temperature	Ratio
C	
25	0.35

Figure 14.33 shows the “PSD” tab window where the outlet PSD is defined in terms of a built-in known distribution function; namely, the normal distribution.

PSD calculation option

- ☐ Copy PSD from inlet
- ☐ Calculate PSD from growth kinetics
- ☐ Calculate PSD from particle growth model
- ☒ User-specified PSD
 - ☐ Overall
 - ☒ Substream ID: **CIPSD**

Crystallizer volume: 4 cum

User-specified PSD

- ☒ Use distribution function
- ☐ User-specified PSD

Bypass fraction: 0

Distribution function ID: 1

Distribution function: Normal

Select parameters: D50

D50: 2 cm



Results

Summary	Balance	Profiles	PSD Results	Status
Crystallizer temperature	101.937	C		
Heater duty	300	kW		
Net duty	0	Gcal/hr		
Crystallizer pressure	1	bar		
Crystallizer volume	4	cum		
Residence time	7.80996	hr		
Crystal product	133.127	kg/hr		
Vapor flow rate	377.506	kg/hr		
Recirculation flow rate	0	kg/hr		
Magma density	259.928	gm/l		

Material	Heat	Load	Vol.% Curves	Wt. % Curves	Petroleum	Polymers	Solids
		Units	FEED	PRODUCT	VAPOR		
▶ Temperature	C	25	101.937	101.937			
▶ Pressure	bar	1	1	1			
▶ Molar Vapor Fraction		0	0	1			
▶ Molar Liquid Fraction		1	0.924363	0			
▶ Molar Solid Fraction		0	0.0756368	0			
▶ Mass Vapor Fraction		0	0	1			
▶ Mass Liquid Fraction		1	0.78614	0			
▶ Mass Solid Fraction		0	0.21386	0			
▶ Molar Enthalpy	kcal/mol	-70.7673	-71.9398	-57.1363			
▶ Mass Enthalpy	kcal/kg	-3153.66	-2728.42	-3171.55			
▶ Molar Entropy	cal/mol-K	-37.5682	-32.1967	-8.73304			
▶ Mass Entropy	cal/gm-K	-1.67418	-1.22111	-0.484757			
▶ Molar Density	mol/cc	0.0509796	0.0460964	3.20657e-05			
▶ Mass Density	kg/cum	1143.97	1215.41	0.577673			



Appendix

Unit operations models used to describe treatment steps carried out on solids in the form of formation, size modification, size-based separation, washing, drying, and fluidization. Aspen Plus built-in unit operation solid models are

1. Crystallizer: produces crystals from solution based on solubility.
2. Crusher: breaks down solid particles to a smaller size.
3. Screen: separates solid particle based on their particle size.
4. Swash: separates solid particles from an entrained liquid of a solids stream using a washing liquid.
5. CCD: separates solid particles from an entrained liquid of a solids stream using a washing liquid in a countercurrent decanter or a multistage washer.
6. Dryer: evaporates volatile moisture components from wet solids.
7. Spray Dryer: evaporates moisture to form particles from sprinkled droplets.
8. Granulator: increases the size of solid particles.
9. Classifier: separates solid particles based on settling velocity.
10. Fluidized Bed: considers both chemical reactions and fluid mechanics.



Solids Separators Models

Unit operation models for separating solids from gases and/or liquids are shown in Table 14.A.1.

TABLE 14.A.1 Unit Operation Models for Solid/Liquid and/or Solid/Gas Separation.

Model	Description	Purpose	Use For
Cyclone	Cyclone separator	Separates solids from gas using gas vortex in a cyclone	Rating and sizing cyclones
VScrub	Venturi scrubber	Separates solids from gas by direct contact with an atomized liquid	Rating and sizing venturi scrubbers
CFuge	Centrifuge filter	Separates solids from liquid using a rotating basket	Rating or sizing centrifuges
Filter	Rotary vacuum filter	Separates solids from liquid using a continuous rotary vacuum filter	Rating or sizing rotary vacuum filters
HyCyc	Hydrocyclone	Separates solids from liquid using liquid vortex in a hydrocyclone	Rating or sizing hydrocyclones
FabFl	Fabric filter	Separates solids from gas using fabric filter baghouses	Rating and sizing baghouses
ESP	Electrostatic precipitator	Separates solids from gas using an electric charge between two plates	Rating and sizing dry electrostatic precipitators

Solids Handling Models

Pneumatic conveyance of granular solid materials over short and long distances through pipes. This includes solids transport through a single pipe or pipeline network.



SOLIDS CLASSIFICATION

Solids are classified by Aspen Plus as shown in Table 14.B.1.

TABLE 14.B.1 Solids Classification Based on Knowledge of Molecular Structure and Chemical Reactivity.

Class	Type	Characteristics	Example
Conventional (a solid with a well-defined molecular structure)	Salts	Participates in phase equilibrium thus defined through chemistry	NaCl(s), ice, and purified terephthalic acid (pTA(s))
	Conventional Inert Solids (CISOLIDS)	An inert solid phase and does not participate in phase equilibrium	SiO ₂ (s) and urea(s)
Non-conventional (complex structure)	Non-conventional Solids (NCSOLIDS)	Characterized through component attributes (ultimate analysis) - special thermodynamic models	Coal and paper pulp

Figure 14.B.1 shows how one can define the same component using different types; hence, a different “Component ID” will be created each time. For example, if “NaCl” is defined as “*Conventional*”, then it will be part of the aqueous medium (i.e., participates in phase equilibrium); on the other hand, if it is defined as “NaCl(S)”, that is, “*Solid*”, then it will be treated as inert (i.e., will not be part of the aqueous medium). Silica being defined as “*Solid*” means it is with a known molecular structure; on the other hand, coal being defined as “*Nonconventional*” means that it has a complex structure.

PREDEFINED STREAM CLASSIFICATION

Selection

Petroleum

Nonconventional

Enterprise Database

Comments

Select components

Component ID	Type	Component name	Alias	CAS number
WATER	Conventional	WATER	H2O	7732-18-5
NACL	Solid	SODIUM-CHLORIDE	NACL	7647-14-5
NA+	Conventional	NA+	NA+	
CL-	Conventional	CL-	CL-	
PTA	Conventional	TEREPHTHALIC-ACID	C8H6O4-D3	100-21-0
PTA(S)	Conventional	TEREPHTHALIC-ACID	C8H6O4-D3	100-21-0
COAL	Nonconventional			
SIO2	Solid	SILICON-DIOXIDE	SIO2	14808-60-7
★				

Find

Elec Wizard

SFE Assistant

User Defined

Reorder

Review

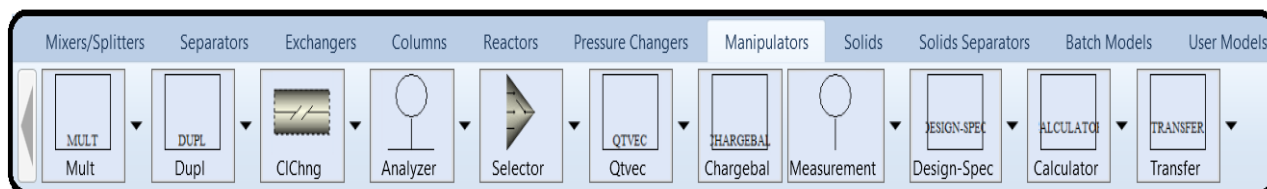


Stream classes are very useful when modeling solids because they can be used to differentiate between the properties of each substream in the simulation. Stream classes ease the integration of solids and fluids in one simulation. The stream class can either be created or selected from a predefined type. Upon one's need, a predefined class can also be modified. Aspen Plus predefined stream classes are in general sufficient for most applications. All unit operation models, except "Extract", can handle stream classes with solid substreams. Table 14.C.1 shows different Aspen Plus built-in stream classes.

TABLE 14.C.1 Aspen Plus Built-In Stream Classes.

Stream Class	When
CONVEN	The simulation does not involve solids, or the only solids are electrolytes salts
MIXCISLD	Conventional inert solids are present, but there is no particle size distribution
MIXNC	Non-conventional solids are present, but there is no particle size distribution
MIXCINC	Both conventional inert and non-conventional solids are present, but there is no particle size distribution
MIXCIPSD	Conventional inert solids are present, with a particle size distribution
MIXNCPSD	Non-conventional solids are present, with a particle size distribution

The stream class can be changed to a different class using the stream class changer (manipulator) as shown in Figure 14.C.1.



SUBSTREAM CLASSES

Substream classes are the building blocks of stream classes. They can either be predefined or customized. Predefined substream classes are shown in Table 14.D.1:



TABLE 14.D.1 Substream Classes for Solid-Bearing Streams.

#	Substream Class	Description
1	MIXED	Fluids only
2	CISOLID	Conventional inert solids that participate in reactions but do not participate in phase equilibria (except in the RGibbs model)
3	NC	Non-conventional solids
4	NCPSD	Non-conventional solids with PSD
5	CIPSD	Conventional inert solids with PSD

Notice that PSD can be specified for each substream in a stream class. Such an advantage permits the user to have different PSD functions for different streams or flowsheet sections. On the other hand, customized substream classes are additional substream types that can be created by the user to complement the existing predefined substreams. Customized substreams are usually used to ease the treatment of different solids in one simulation and permit the use of different PSDs for one predefined substream type, as shown in Figure 14.D.1.

PARTICLE SIZE DISTRIBUTION (PSD)

In Aspen Plus, particle size distribution (PSD) is represented by the weight fractions per particle size interval, given the number of intervals and the size range for each interval. In other words, a PSD describes the amount of particles in a sample of material with respect to size. The built in Aspen Plus particle size distribution has 10 predefined size intervals. The user can modify the built in particle size distribution by changing the number of intervals or the size ranges for the intervals. In some situations, he/she may want to have two or more particle size distribution definitions, with different size ranges. This will be useful if different sections of the flowsheet have different particle sizes. In this regard, a PSD mesh is defined (i.e., a grid with size intervals over which the particle sizes will be described). The PSD is populated via experimental results (i.e., user-specified weight fractions) or utilizing built-in distribution functions. In Aspen Plus, there are simulation PSD meshes coined for a substream under which all the results for multiple streams will be presented and input stream PSD meshes that are defined for individual feed streams. For user-specified PSD meshes, Aspen Plus provides four types: equidistant, geometric, logarithmic, and user-defined, where the user is required to enter the lower limit, upper limit, and size unit (i.e., mm, μm , angstrom, etc.). Depending on the selected type, the user may also have to enter the number of intervals. For the user-defined type, it is suited for experimental data where the user can copy and paste values from other spreadsheet packages such as Microsoft Excel sheet. For built-in distribution functions, the user, however, must specify the function parameters that describe, in general, the mean or a reference datum and the degree of scatter around that mean or reference value, as in the following distributions:



1. Gates–Gaudin–Schuhmann (GSS): It has two parameters: one for the maximum particle size (i.e., at which the cumulative mass fraction $Q(d)=1$) and another for describing the profile (shape) of the distribution itself (i.e., narrow vs. broad). GGS is suited for coarse grinding.

2. Rosin–Rammler–Sperling–Bennet (RRSB): It has two parameters: one for the particle size (i.e., at which the cumulative mass fraction $Q(d)=0.632$) and another for describing the profile of the distribution itself (i.e., narrow vs. broad). RRSB is suited for fine grinding.

3. Normal: It has two parameters: one for the mean or median (at which the cumulative mass fraction $Q(d)=0.5$) and another for describing the profile of the distribution itself (i.e., narrow vs. broad).

4. Log Normal: It has two parameters: one for the median (the median at which the cumulative mass fraction $Q(d)=0.5$) and another for describing the profile of the distribution itself (i.e., narrow vs. broad). If particles are broken up at random, they will end up with such a distribution. When viewing and analyzing the results for solid components, Aspen Plus provides sets of properties:

1. Median value (D50): the size at which 50% of particles are larger.

2. Mean particle size (D1_0): the sum over all intervals for the multiplication of the average particle size, in one interval, by the fraction.

3. Specific surface area (VSSA): the ratio of the surface area of a particle to its volume.

For a completely spherical object, it is $4\pi r^2 / \frac{4}{3}\pi r^3 = \frac{3}{r} = \frac{6}{D}$. For other non-spherical objects, a shape or sphericity factor is used.

4. Sauter mean diameter (D3_2 or SMD): an average particle size. SMD is the diameter of spheres with the same volume/area ratio as the particle mixture. For example, if V is the total volume of particles and A is their total surface area, then $SMD=6V/A=D_{particle}$. SMD is inversely proportional to VSSA. SMD is the characteristic diameter for a packed bed flow and is mainly used in fluidized bed calculations.

FLUIDIZED BEDS

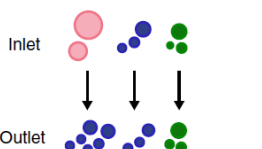
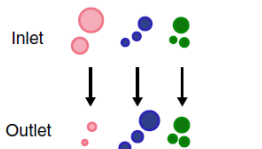
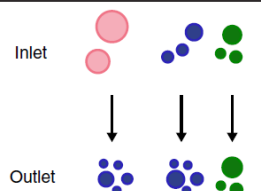
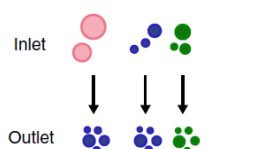



The Aspen Plus fluidized bed model describes a bubbling or circulating fluidized bed and tackles five different aspects: Entrainment of Particles: It takes into account the geometry of the vessel and additional gas supply and provides options to determine the minimum fluidization velocity, transport disengagement height, and distributor pressure drop for either a porous plate or bubble cap.

Chemical Reactions: After defining the reaction stoichiometry and its kinetics, the calculation method treats the gas as plug flow, the solids ideally mixed, and each balance cell is considered as CSTR (i.e., uniform in properties).

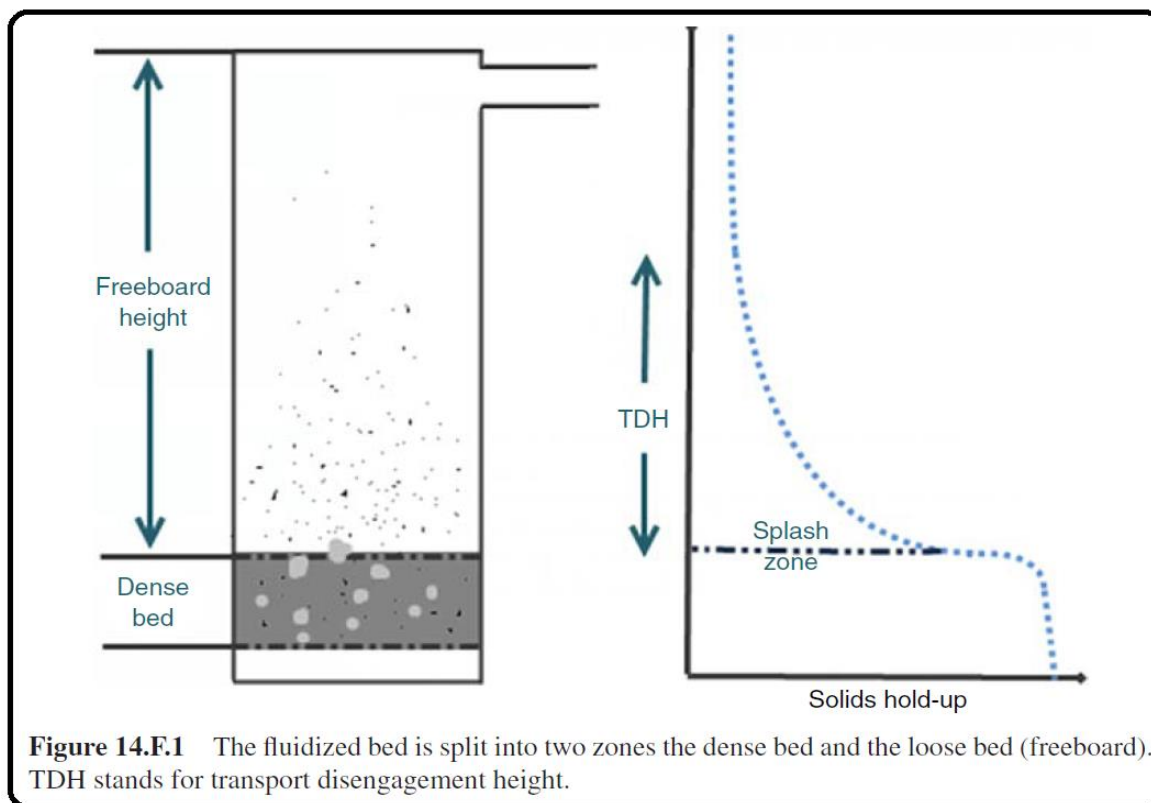
Fluid Mechanics: It assumes one-dimensional fluidmechanics and considers the impact of volume production/or reduction and heat exchange on fluid mechanics. Change in PSD: Table 14.F.1 demonstrates different options of tackling the outlet PSD with respect to the inlet PSD.



Thermodynamics: Solids and vapors are in thermodynamic equilibrium while the model considers the impact of heat exchange and heat effect on bed temperature.

TABLE 14.F.1 Different Approaches in Tackling the Outlet PSD For a Fluidized Bed.			
Keep PSD	Constant # of Particles	Overall PSD	PSD Per Particle Type
 <p>Particle types have the same PSD as the inlet and the new generated particle types have the overall PSD from the inlet. Total number of particles may change and inert particles will keep the same size</p>	 <p>Total number of particles is fixed. The particle may shrink or grow, depending on the mass loss/gain of the particle type. Inert particles will keep the same size</p>	 <p>Overall bed PSD is defined by the user. Total number of particles may change. Inert particles are allowed to change their size</p>	 <p>Bed PSD is defined by the user on a particle-type basis. Total number of particles may change. Inert particles will keep the same size if PSD is defined only for the reacting species</p>
<p>Reaction: A  \rightarrow B </p> <p>Inert: C </p>			

The Aspen Plus fluidized bed model treats the bed as made of two zones. See Figure 14.F.1.





The bottom zone (dense bed): High solids concentration and fluid mechanics is dealt with according to Werther and Wein [1], which considers the growth and splitting of bubbles. Bubble-related profiles (e.g., bubble diameter, bubble rise velocity, interstitial gas velocity, pressure, and solid volume concentration profiles) can all be calculated.

The freeboard zone (loose bed): Relatively low solids concentration and fluid mechanics is dealt with according to Kunii and Levenspiel [2]. Using the selected entrainment correlation, the solids mass flow and PSD at the outlet condition can be calculated. Overall, once the user defines bed inventory, by specifying the pressure drop or the solids hold-up, the height of the bottom zone and the freeboard can be determined.



Reference

1. Our team experience
2. Werther, J. and Wein, J. (1994) Expansion behavior of gas fluidized beds in the turbulent regime. AIChE Symposium Series, **90** (301), 31–44.
3. Kunii, D. and Levenspiel, O. (1991) Fluidization Engineering, Butterworth-Heinemann, Boston.
4. Aspen Plus – Chemical Engineering Application by KAMAL I.M. AL-MALAH