# ADMS 6 Complex Terrain Validation Clifty Creek Power Plant

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April 2023

### 1 Introduction

The Clifty Creek power plant<sup>1</sup> is located in southern Indiana on the north side of the Ohio River (see **Figure 1**). The area immediately north of the facility is characterized by cliffs rising about 115 m above the river and intersected by creek valleys. The tops of the stacks extend about 80 to 100 m above the top of the cliffs on both sides of the river. Three 208 m stacks were modeled in this evaluation.

There were six SO<sub>2</sub> monitors on the surrounding terrain. One was located in the river valley approximately 8 km to the east (upriver) at about the same elevation as the sources. Another was located 3 km south of the facility on Liberty Ridge near the meteorological tower on the south side of the Ohio River and about 110 m above stack base. The remaining four monitors were located on the terrain north and north-east of the facility at about 125 m above the base of the stacks, ranging in distance from four to 15 km.

Meteorological data from this field study covered 1 January 1975 through to 31 December 1975. The on-site meteorological data were recorded on an instrumented meteorological tower 3 km south of the facility (across the river in Kentucky) on Liberty Ridge.

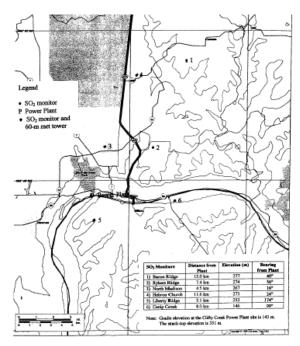


Figure 1 – Monitoring locations in the vicinity of the Clifty Creek power plant.

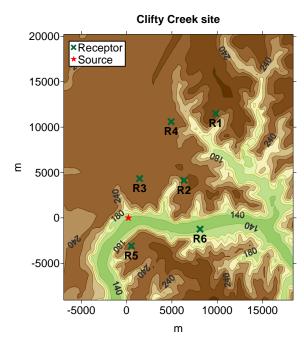


Figure 2 – Modelled area around the Clifty Creek power plant (met. tower close to R5 monitor.)

<sup>&</sup>lt;sup>1</sup> Note that the study description and **Figure 1** have been taken directly from the document [1].



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The input data for the ADMS runs were taken from the AERMOD files downloaded from the United States Environmental Protection Agency website [2]. These data include the observed concentrations that are used for comparison with the ADMS modelled concentrations.

This document compares the results of ADMS 5.2.0.0 (hereafter referred to as ADMS 5.2) with those of ADMS 6.0.0.1 (hereafter referred to as ADMS 6.0).

Section 2 describes the input data used for the model. The results are presented in Section 3 and discussed in Section 4.

## 2 Input data

#### 2.1 Study area

The study area is located around 38.7°N. The roughness length varies, depending on wind direction and month as shown in **Table 1**.

Terrain data included in the modelling covered a  $20 \text{ km} \times 28 \text{ km}$  area (**Figure 2**) with terrain data points located every 300-400 m.

Time of the year	Wind sector				
Time of the year	245-185°	185-245°			
January to March	0.20	0.40			
April to May	0.30	0.50			
June to August	0.40	0.60			
September to November	0.30	0.50			
December	0.20	0.40			

**Table 1** – Surface roughness length (m) according to wind sector.

#### 2.2 Source parameters

The source parameters are summarised in **Table 2**. The minimum exit velocity for all stacks is just under 25 m/s. The maximum exit velocity for stack 1 is 57 m/s, whereas stacks 2 and 3 have exit velocity peaks of 103.5 and 102.7 m/s. Emissions rates vary from 3-3682 g/s.

Source name	Pollutant	Location	Stack height (m)		Exit T (°C)	Diameter (m)	Emission rate (g/s)
Stack 1	$SO_2$	200, 0	207.9	varied	172.22	4.63	varied
Stack 2	$SO_2$	200, 0	207.9	varied	172.22	4.63	varied
Stack 3	$SO_2$	200, 0	207.9	varied	172.22	4.63	varied

**Table 2** – Source input parameters. T is the temperature, V the velocity.

#### 2.3 Receptors

The receptor network consists of 6 monitors, ranging from 4 to 15 km from the sources. All receptors are modelled as ground level receptors.

Figure 2 shows the receptor network used in the experiment. Table 3 gives the receptor



coordinates in a local coordinate system; for reference, the stacks are located at (200,0).

<b>Receptor name</b>	Location
R1	9842, 11491
R2	6335, 4138
R3	1440, 4326
R4	4918, 10597
R5	524, -3083
R6	8102, -1251

Table 3 – Receptor point locations.

#### 2.4 Meteorological data

One year of hourly sequential data from 1 January 1975 to 31 December 1975 was used. The wind rose is shown in **Figure 3**. **Table 4** gives the detail of the modelled conditions. The criteria for the stability categories are as follows, where H is the boundary layer height and  $L_{MO}$  is the Monin-Obukhov length, as calculated by the model's meteorological processor:

 $Stable: H/L_{MO} > 1 \\ Neutral: -0.3 \leq H/L_{MO} \leq 1 \\ Convective: H/L_{MO} < -0.3 \\$ 

	Conditions	<b>ADMS 5.2</b>	<b>ADMS 6.0</b>
Hours modelled	Stable conditions	5012 (59%)	5032 (59%)
	Neutral conditions	943 (11%)	918 (11%)
	Unstable conditions	2556 (30%)	2566 (30%)
	Total	8511 (100%)	8516 (100%)
	Calm conditions	0	0
Hours not	Wind speed at $10 \text{ m} < 0.75 \text{ m/s}$	248	243
modelled	Inadequate data	1	1
	Total	249	244

**Table 4 –** Meteorological conditions. Percentage values are computed with respect to the total number of modelled hours.

The wind speeds vary from 0.27 to 21.1 m/s and the surface temperature from -18.8 to 34.4°C. The height of the recorded wind measurement was 60 m. The wind rose is shown in **Figure 3**.



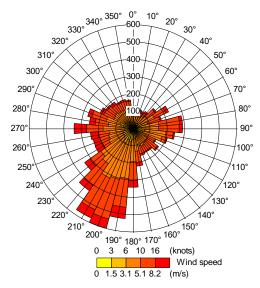


Figure 3 – Wind rose from meteorological data.

#### 3 Results

Scatter plots and quantile-quantile plots of model results against observed data are presented in Section 3.1. Other statistical analysis is presented in Section 3.2. The graphs and statistical analysis have been produced by the Model Evaluation Toolkit v5.2 [4].

#### 3.1 Scatter and quantile-quantile plots

The modelled  $SO_2$  concentrations are compared to observed hourly concentrations ( $\mu g/m^3$ ). **Figures 4** and **5** show frequency scatter plots and quantile-quantile plots of results. Note that these quantile-quantile plots are *linear*; care should be exercised when comparing these plots with similar ones presented with *logarithmic* axes.



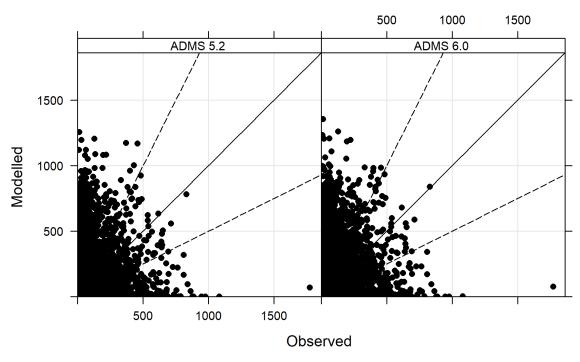


Figure 4 – Scatter plots of ADMS results against observed data for all receptors (µg/m³).

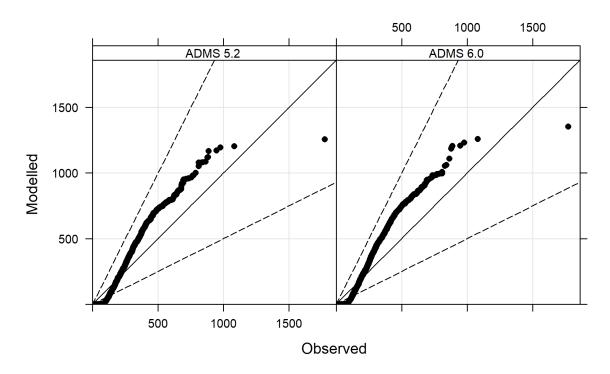


Figure 5 – Quantile-quantile plots of ADMS results against observed data for all receptors (μg/m³).

#### 3.2 Statistics

**Table 5** compares the modelled and observed maximum 1-hour, 3-hour and 24-hour average concentrations at the receptor points. **Table 6** compares the corresponding robust highest

concentrations, where this statistic is defined by:

robust highest concentration = 
$$\chi(n) + (\chi - \chi(n)) \ln \left(\frac{3n-1}{2}\right)$$
,

where n is the number of values used to characterise the upper end of the concentration distribution,  $\chi$  is the average of the n-1 largest values, and  $\chi(n)$  is the  $n^{th}$  largest value; n is taken to be 26, as in Perry et al. [3].

Statistics	stics Data	Concentrations (µg/m³)						Mean M/O
Staustics		<b>P1</b>	P2	P3	P4	P5	P6	ratio
1 have	Observed	887	809	1772	976	692	1081	-
1-hour maximum	ADMS 5.2	720	1204	1195	966	1079	1256	1.11
	ADMS 6.0	801	1259	1185	1009	1053	1355	1.16
2.1	Observed	579	536	857	613	475	627	-
3-hour	ADMS 5.2	567	857	935	734	600	792	1.23
maximum	ADMS 6.0	709	1206	906	700	526	824	1.35
24-hour maximum	Observed	267	195	338	181	153	171	-
	ADMS 5.2	165	159	266	217	155	113	0.85
	ADMS 6.0	194	168	273	219	152	115	0.88

**Table 5** – Observed (O) and modelled (M) maximum concentrations (ug/m³) per receptor point, and the mean ratio of modelled/observed values for each statistic.

Statistics	Data	Robust Highest Concentrations (µg/m³)						Mean M/O
Statistics		<b>P1</b>	P2	P3	P4	P5	P6	ratio
1 have	Observed	892	926	1334	973	523	856	-
1-hour RHC	ADMS 5.2	855	1142	1274	1020	1272	1235	1.35
KHC	ADMS 6.0	917	1236	1222	1059	1216	1386	1.39
3-hour	Observed	664	586	776	607	411	579	-
S-nour RHC	ADMS 5.2	590	899	973	733	620	620	1.24
RHC	ADMS 6.0	695	976	1011	701	614	752	1.33
24-hour RHC	Observed	232	194	208	198	146	154	-
	ADMS 5.2	141	178	272	212	121	120	0.92
	ADMS 6.0	150	182	296	210	121	149	0.98

**Table 6** – Observed (O) and modelled (M) robust highest concentrations (RHC) per receptor point, and the mean ratio of modelled/observed RHC for each statistic (number of points = 26).

#### 4 Discussion

The scatter and quantile-quantile plots (**Figures 4** and **5**) show good agreement between modelled and observed concentrations. The scatter plots compare predicted and measured concentrations at a particular location at a particular time, i.e. an (x,t) pairing. The quantile-quantile plots compare the distribution of predicted and measured concentrations during the period having abandoned the (x,t) pairing. Predicting the distribution of concentrations accurately is relevant to calculations for permitting purposes, where the comparison with air quality limits is more important than accurately predicting a time series of concentrations at each location. The latter is a harder task.

The pollutant monitored for this study is SO<sub>2</sub>. There are a number of issues with using SO<sub>2</sub> as



a tracer, which include:

- The detection limits of monitors are usually of the order of  $16 \,\mu g/m^3$ , and concentrations below these are set to one-half of the limit. This leads to considerable inaccuracy when modelled concentrations are low.
- SO<sub>2</sub> is released from other sources. If estimates of these background concentrations are not available, then the model will underestimate concentrations, particularly long-term averages.

Comparisons between modelled and observed annual average concentrations are not presented in this report due to the issues with monitor detection limits and background data.

The predictions of maximum concentrations and robust highest concentrations presented in **Tables 5** and **6** show good model performance considering the complexity of the domain modelled.

The differences between ADMS 5.2 and ADMS 6.0 are generally small. ADMS 6.0 predicts slightly higher maximum concentrations and RHC values than ADMS 5.2 at all receptors, which leads to slightly worse mean modelled-to-observed ratios, apart from the values averaged over 24 hours. There has been a change to the meteorological processor, in which the solar elevation angle is calculated at the middle of the hour rather than the end of it, which is having some effect in daylight hours.

#### 5 References

- [1] Paine, R.J, Lee, R.F, Brode, R, Wilson, R.B, Cimorelli, A.J., Perry, S.G., Weil, J.C., Venkatram, A, and Peters, W., 1998: Model Evaluation Results for AERMOD (draft). United States Environmental Protection Agency.
- [2] United States Environmental Protection Agency website, Model Evaluation Databases. <a href="https://www.epa.gov/scram/air-quality-dispersion-modeling-preferred-and-recommended-models">https://www.epa.gov/scram/air-quality-dispersion-modeling-preferred-and-recommended-models</a>
- [3] Perry, S. G., Cimorelli, A. J., Paine, R.J., Brode, R.W., Weil, J.C., Venkatram, A., Wilson, R.B., Lee, R.F, & Peters, W.D. 2005 AERMOD: A Dispersion Model for Industrial Source Applications. Part II: Model Performance against 17 Field Study Databases. *J. Appl. Met.* 44, pp 694-708.
- [4] Stidworthy A, Carruthers D, Stocker J, Balis D, Katragkou E, and Kukkonen J, 2013: *MyAir Toolkit for Model Evaluation*. 15<sup>th</sup> International Conference on Harmonisation, Madrid, Spain, May 2013.

