

**Supplementary Information for:
Contributions to local and regional-scale formaldehyde concentrations**

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Notes on units of gas-phase mixing ratios: 1 ppb = 1 nmol/mol; 1 ppt = 1 pmol/mol.



Figure S1: Measurement sites used for the evaluation of the forward model. Sites include ambient air monitoring locations from the air basin- and state-wide regulatory agencies.

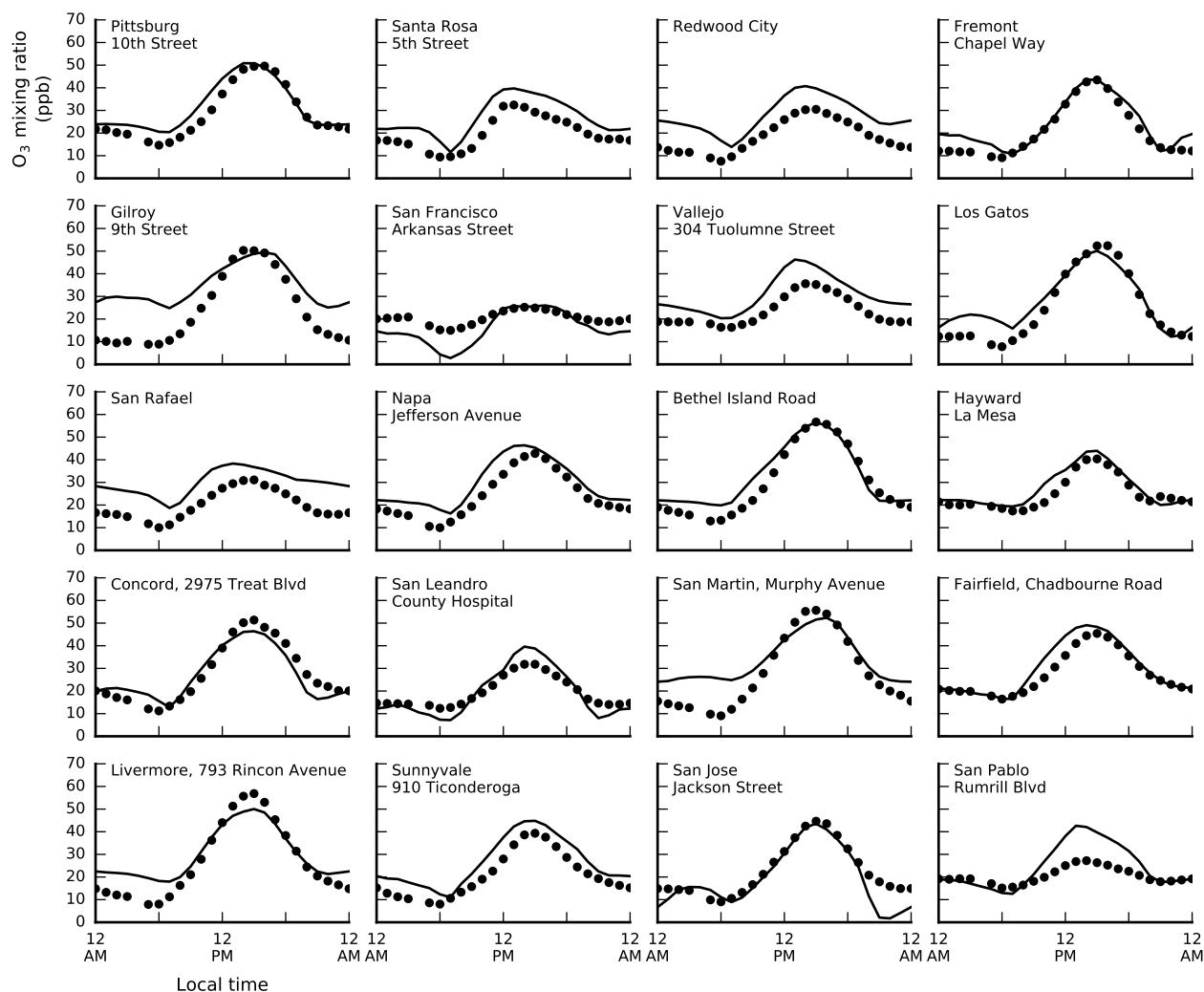


Figure S2: Observed (filled circles) and modeled (solid line) average hourly ozone mixing ratios at twenty different monitoring sites in July. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

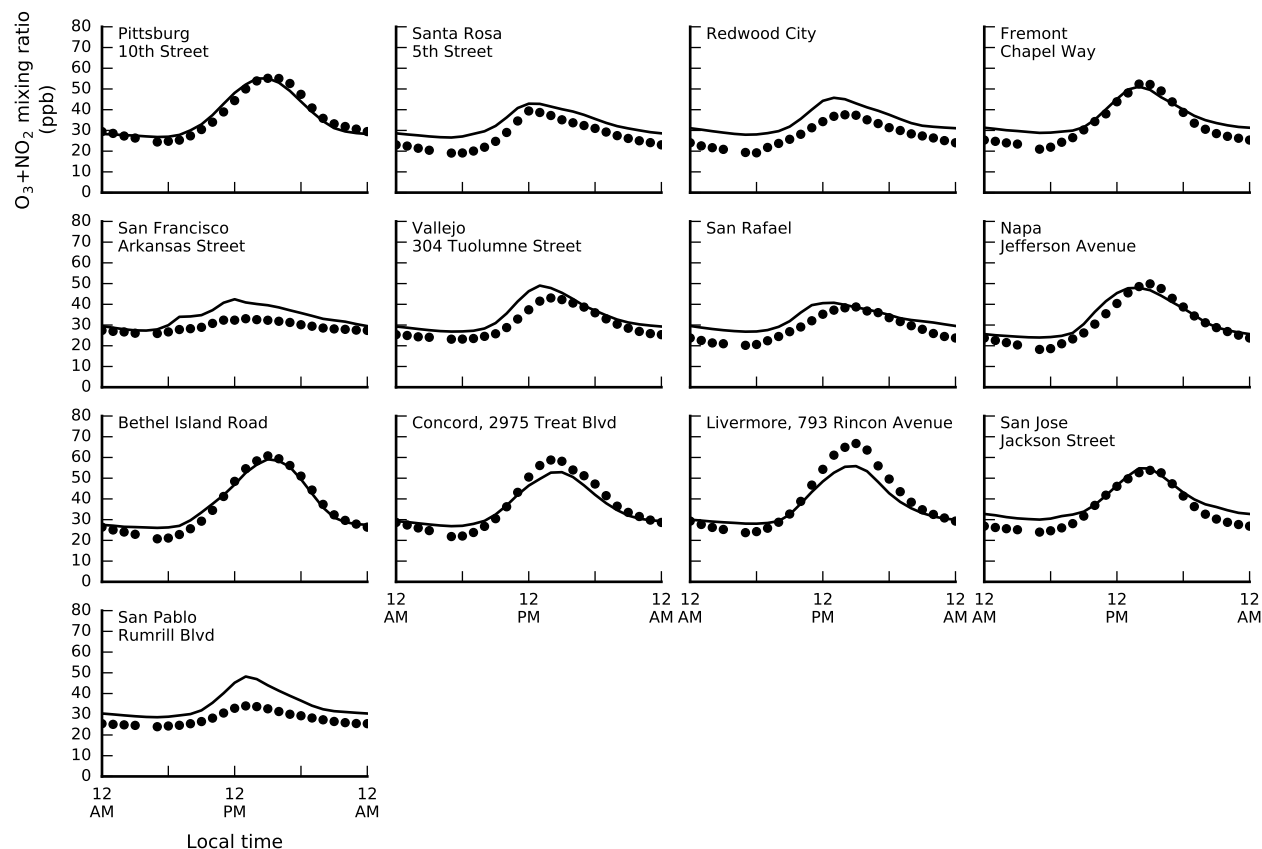


Figure S3: Observed (filled circles) and modeled (solid line) average hourly $O_3 + NO_2$ mixing ratios at thirteen different monitoring sites in July. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

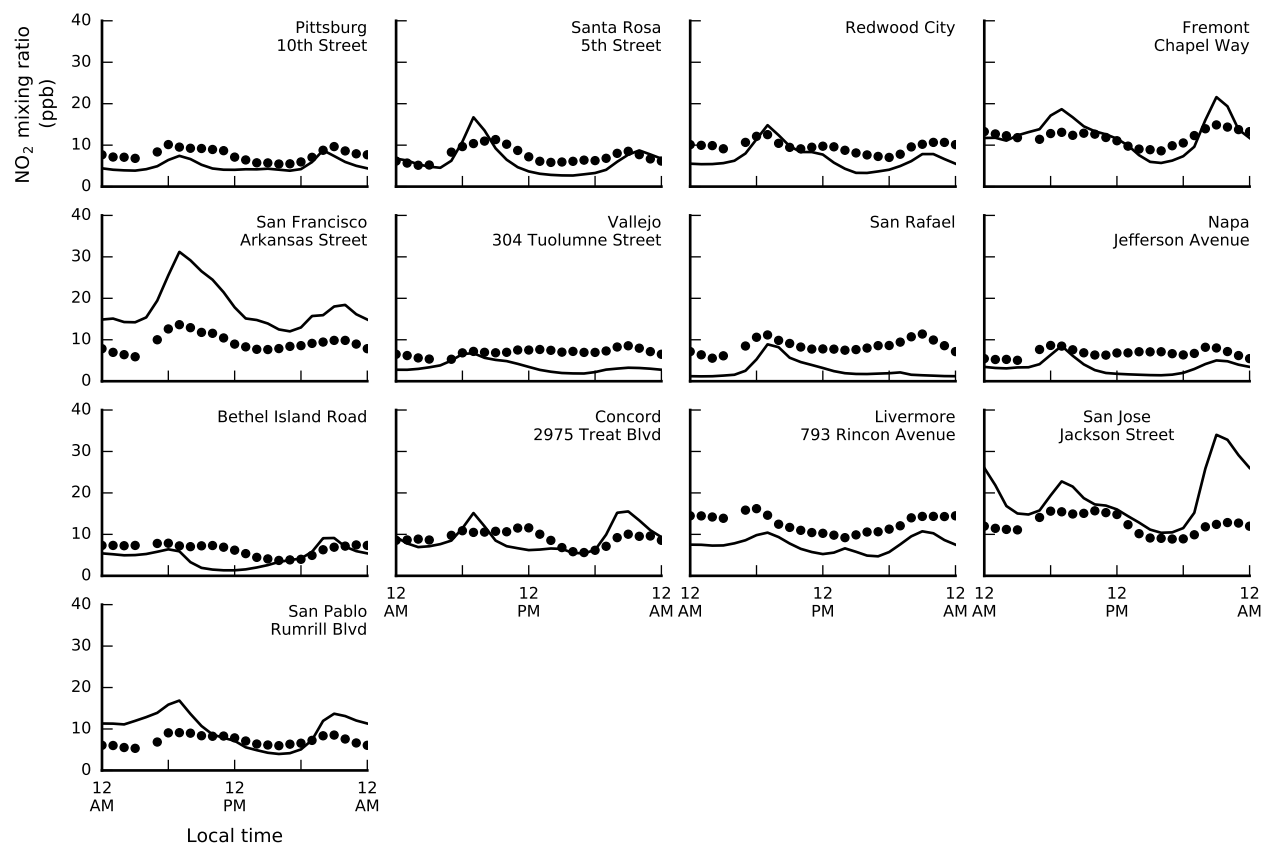


Figure S4: Observed (filled circles) and modeled (solid line) average hourly NO_2 mixing ratios at thirteen different monitoring sites in July. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

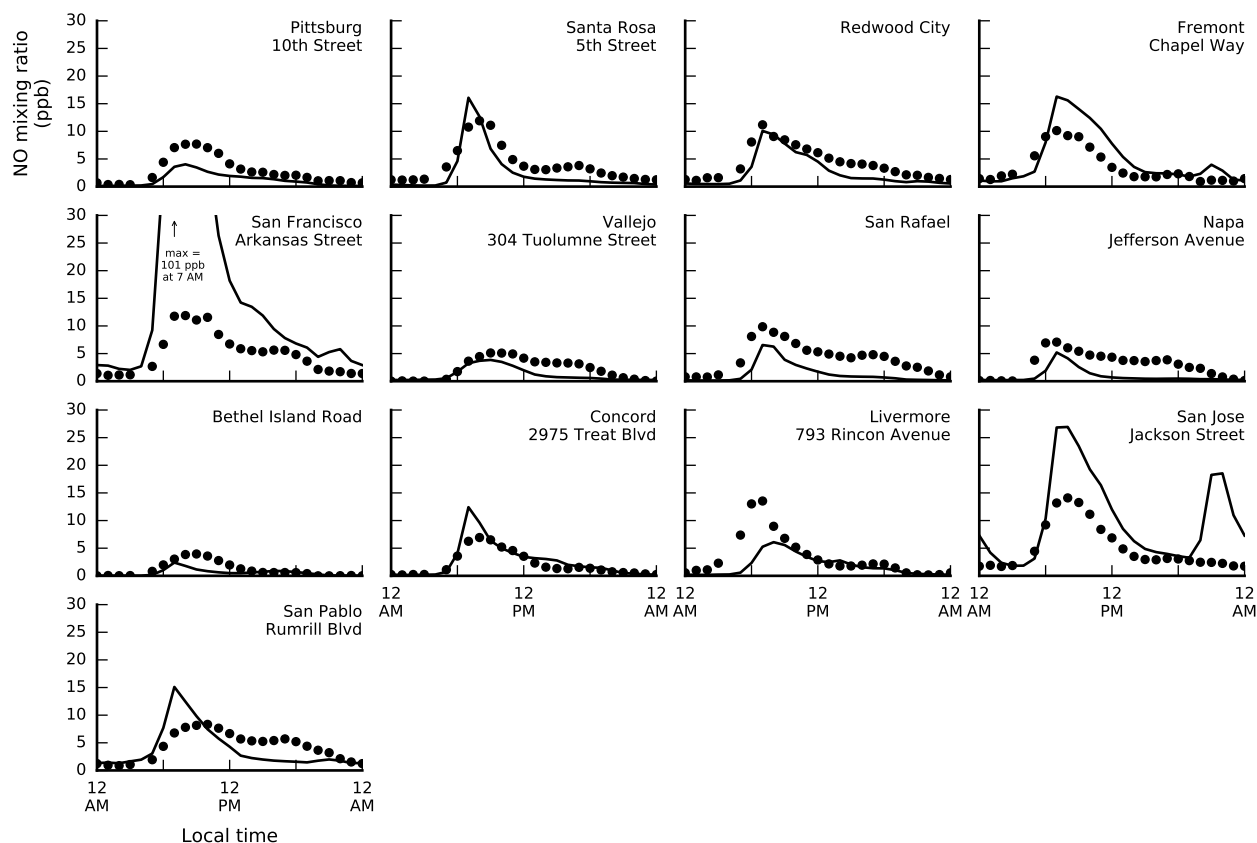


Figure S5: Observed (filled circles) and modeled (solid line) average hourly NO mixing ratios at thirteen different monitoring sites in July. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

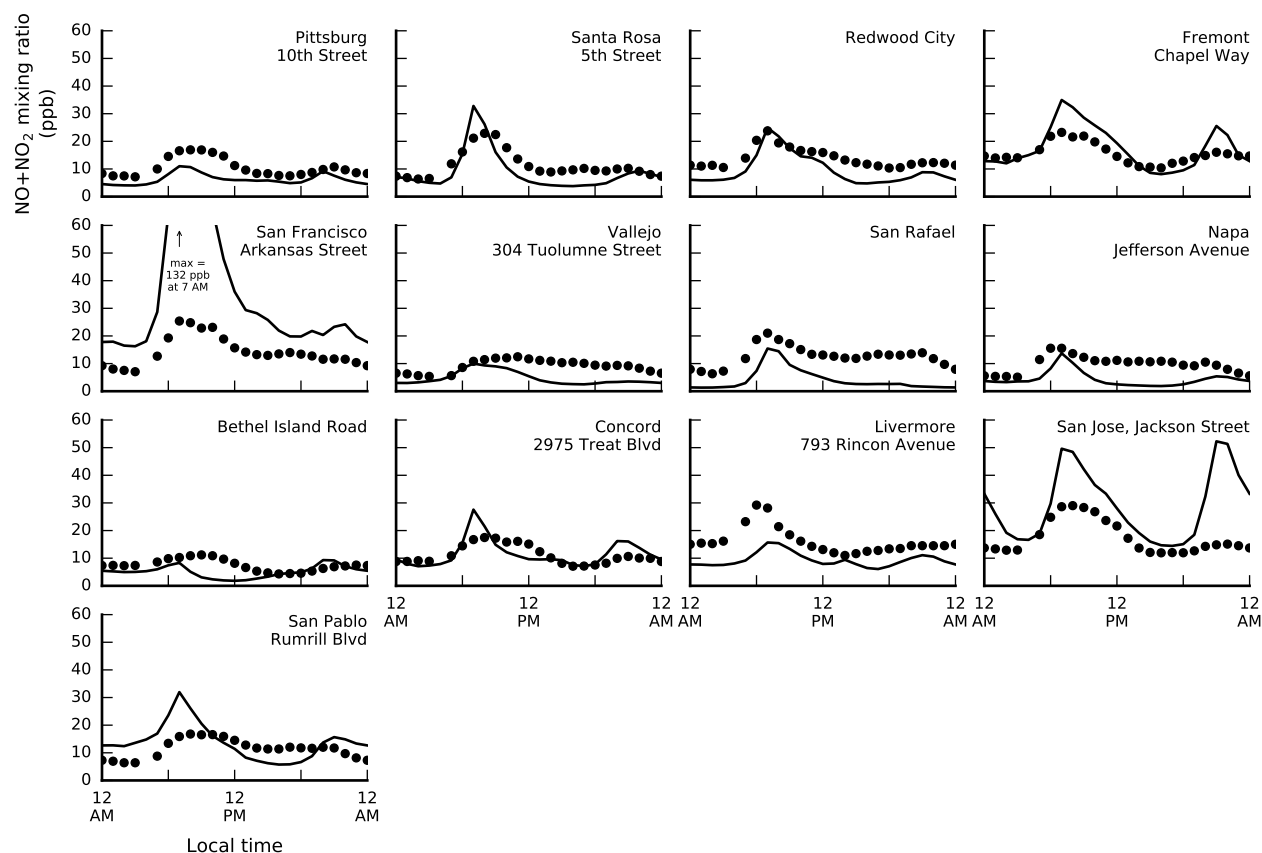


Figure S6: Observed (filled circles) and modeled (solid line) average hourly NO + NO₂ mixing ratios at thirteen different monitoring sites in July. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

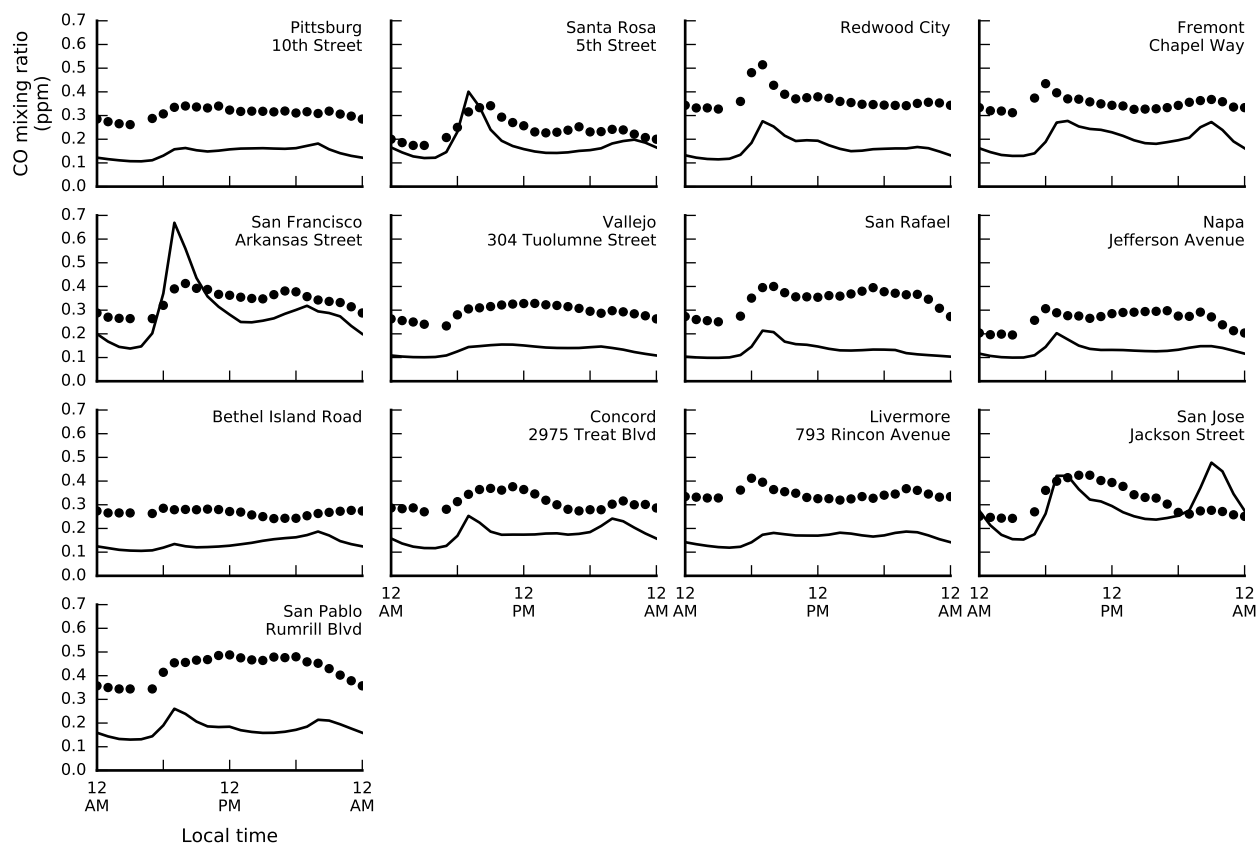


Figure S7: Observed (filled circles) and modeled (solid line) average hourly CO mixing ratios at thirteen different monitoring sites in July. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

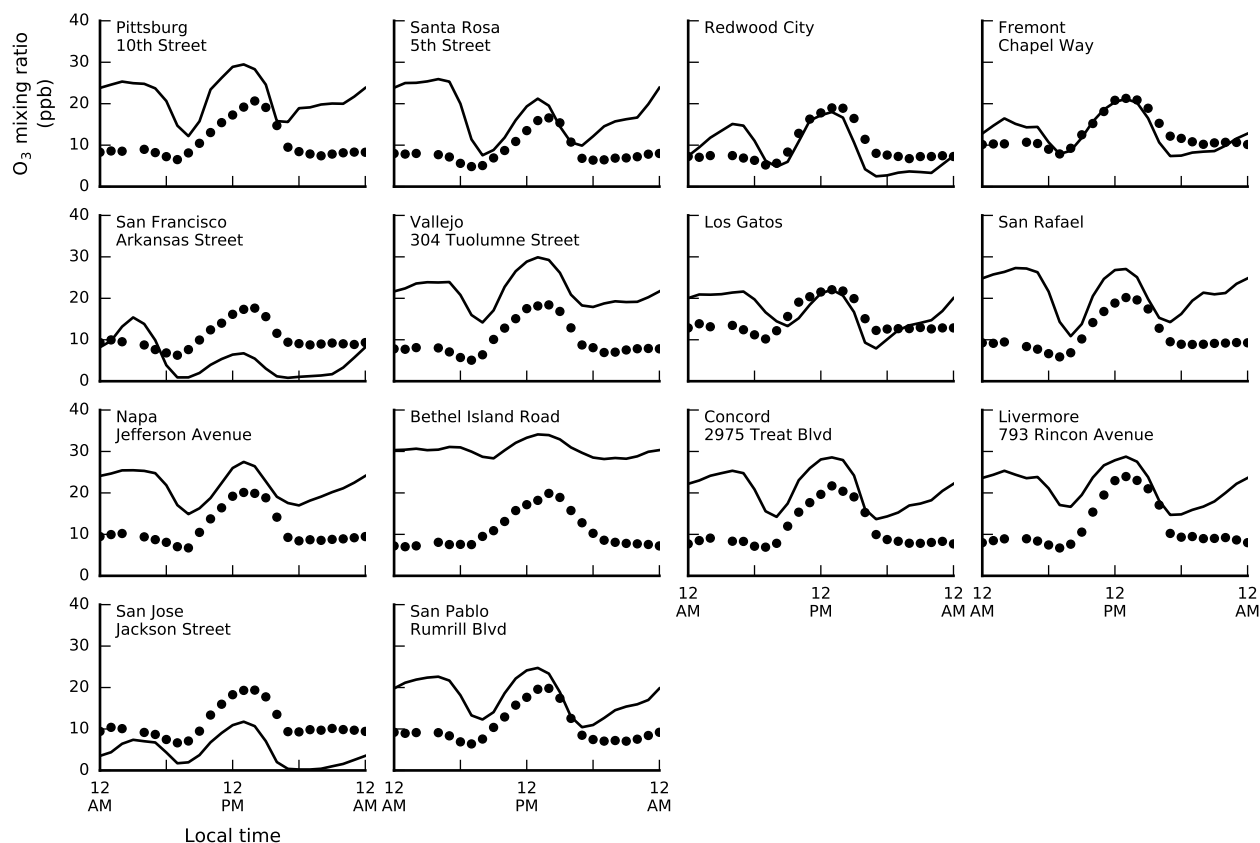


Figure S8: Observed (filled circles) and modeled (solid line) average hourly ozone mixing ratios at fourteen different monitoring sites in December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

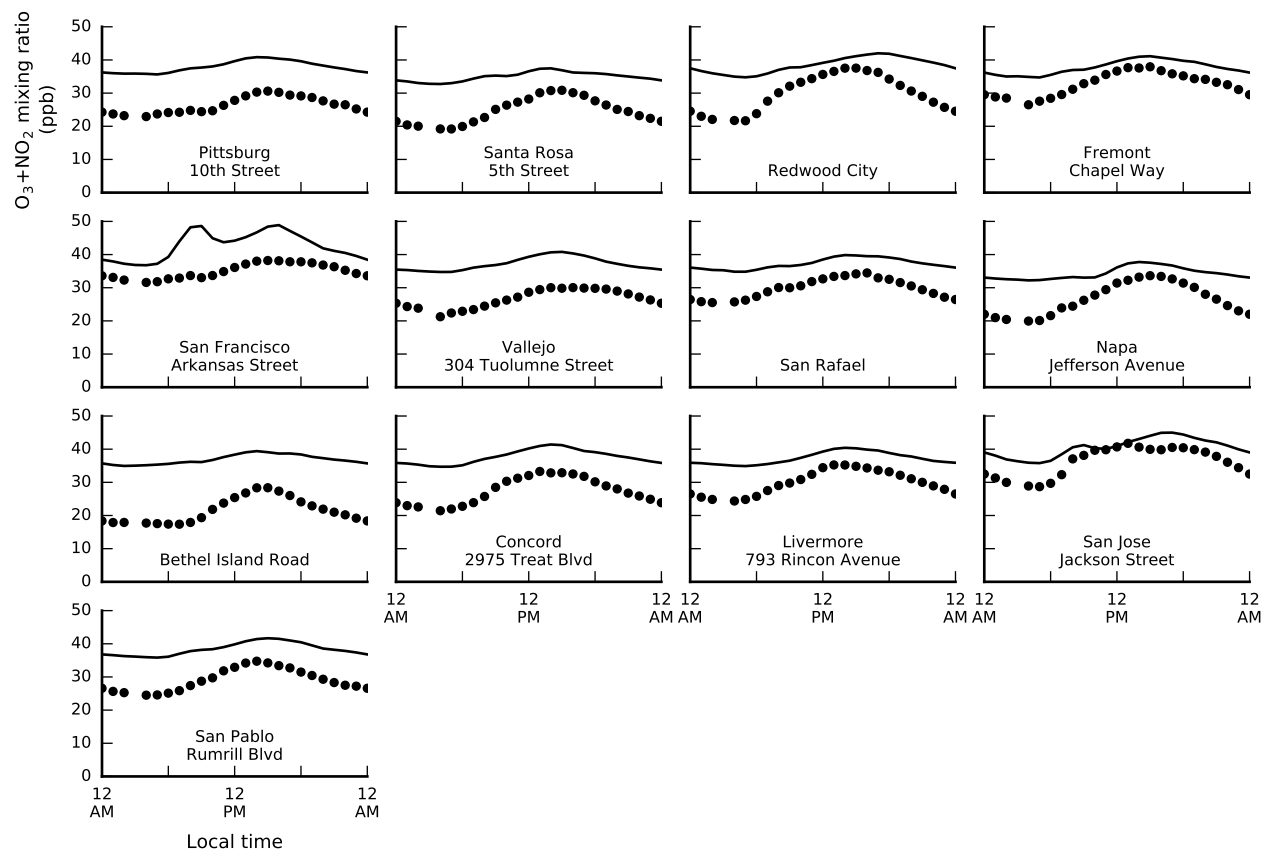


Figure S9: Observed (filled circles) and modeled (solid line) average hourly $O_3 + NO_2$ mixing ratios at thirteen different monitoring sites in December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

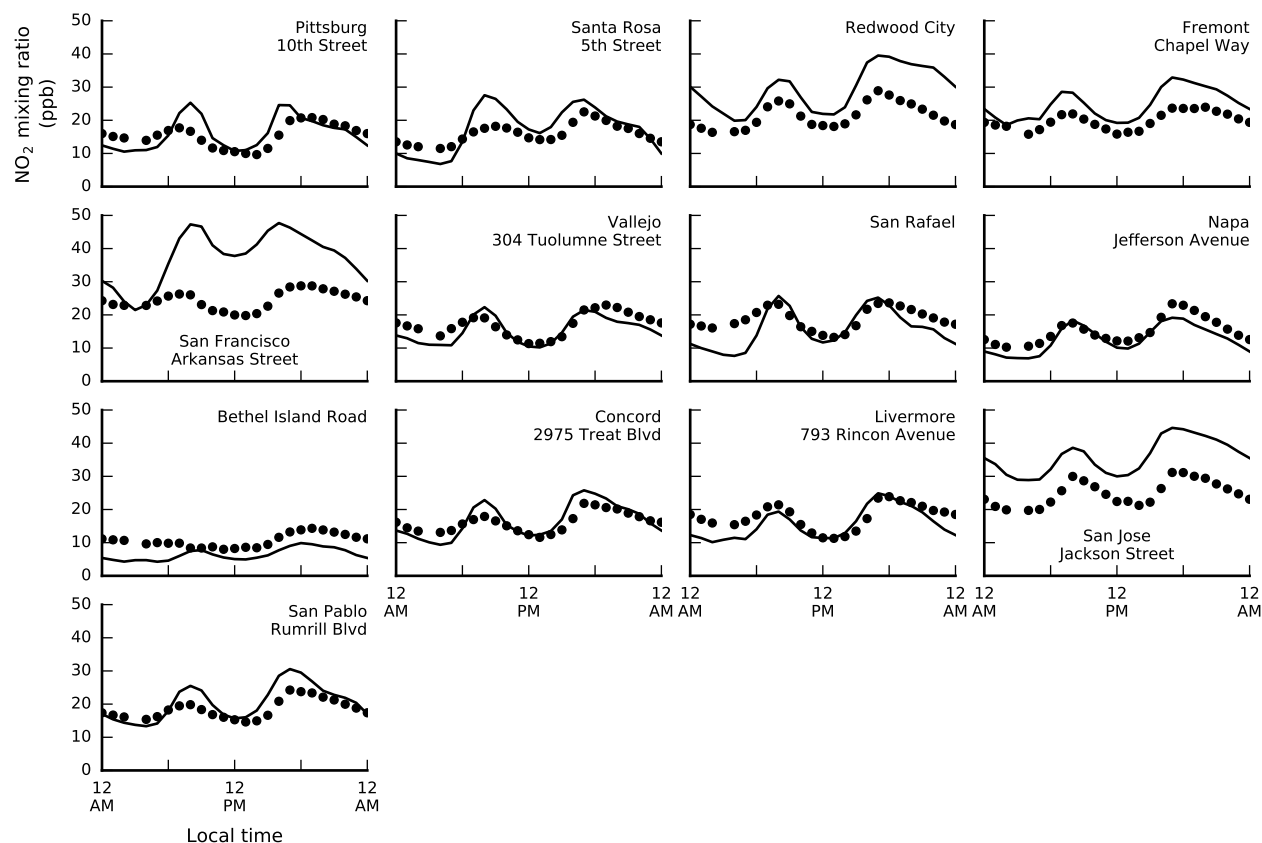


Figure S10: Observed (filled circles) and modeled (solid line) average hourly NO_2 mixing ratios at thirteen different monitoring sites in December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

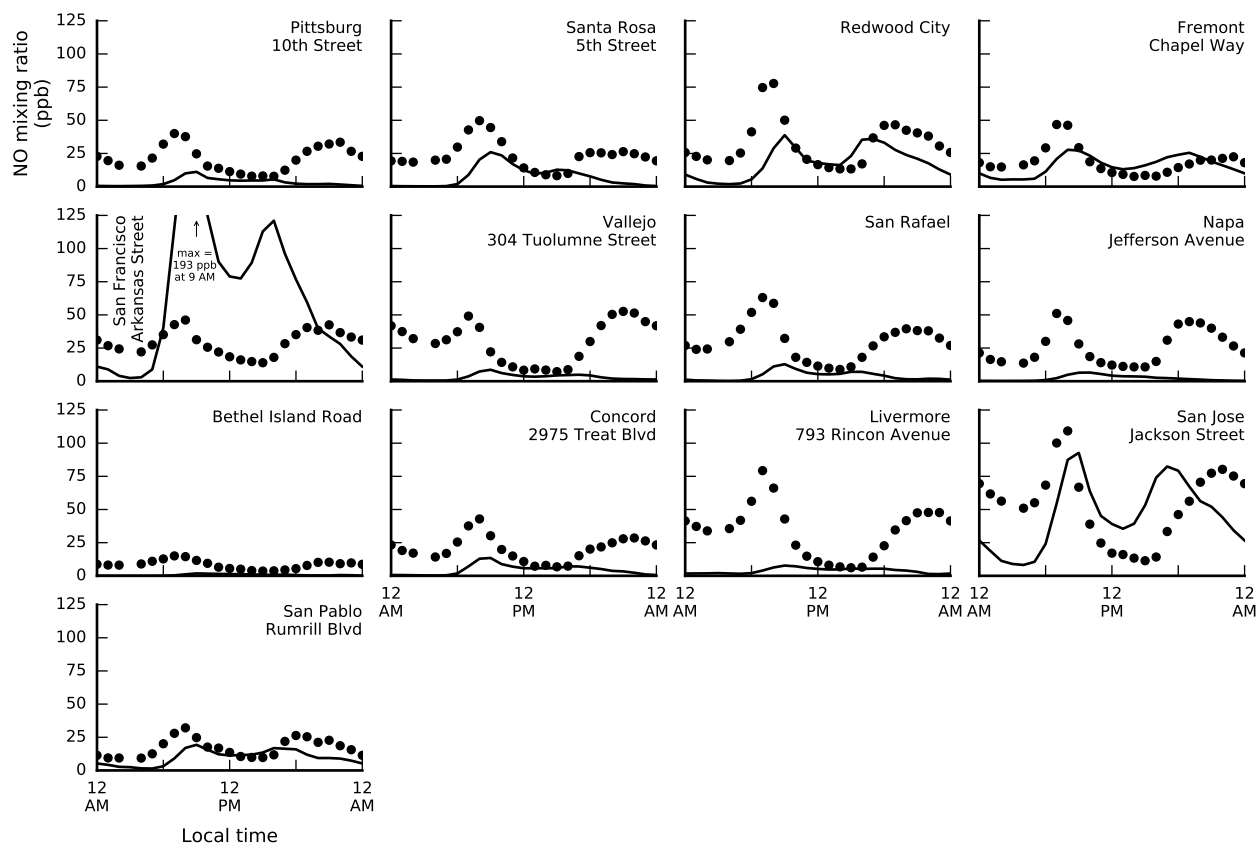


Figure S11: Observed (filled circles) and modeled (solid line) average hourly NO mixing ratios at thirteen different monitoring sites in December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

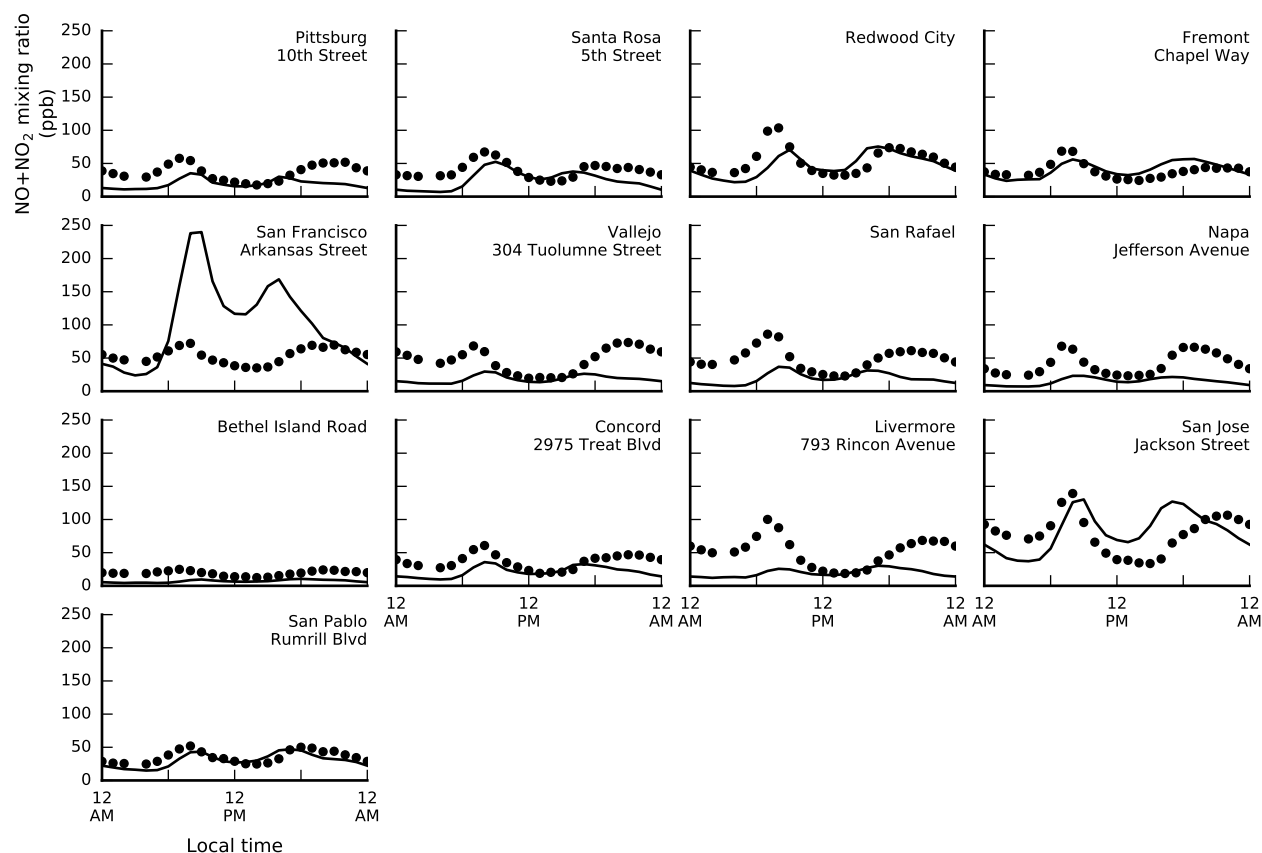


Figure S12: Observed (filled circles) and modeled (solid line) average hourly NO + NO₂ mixing ratios at thirteen different monitoring sites in December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

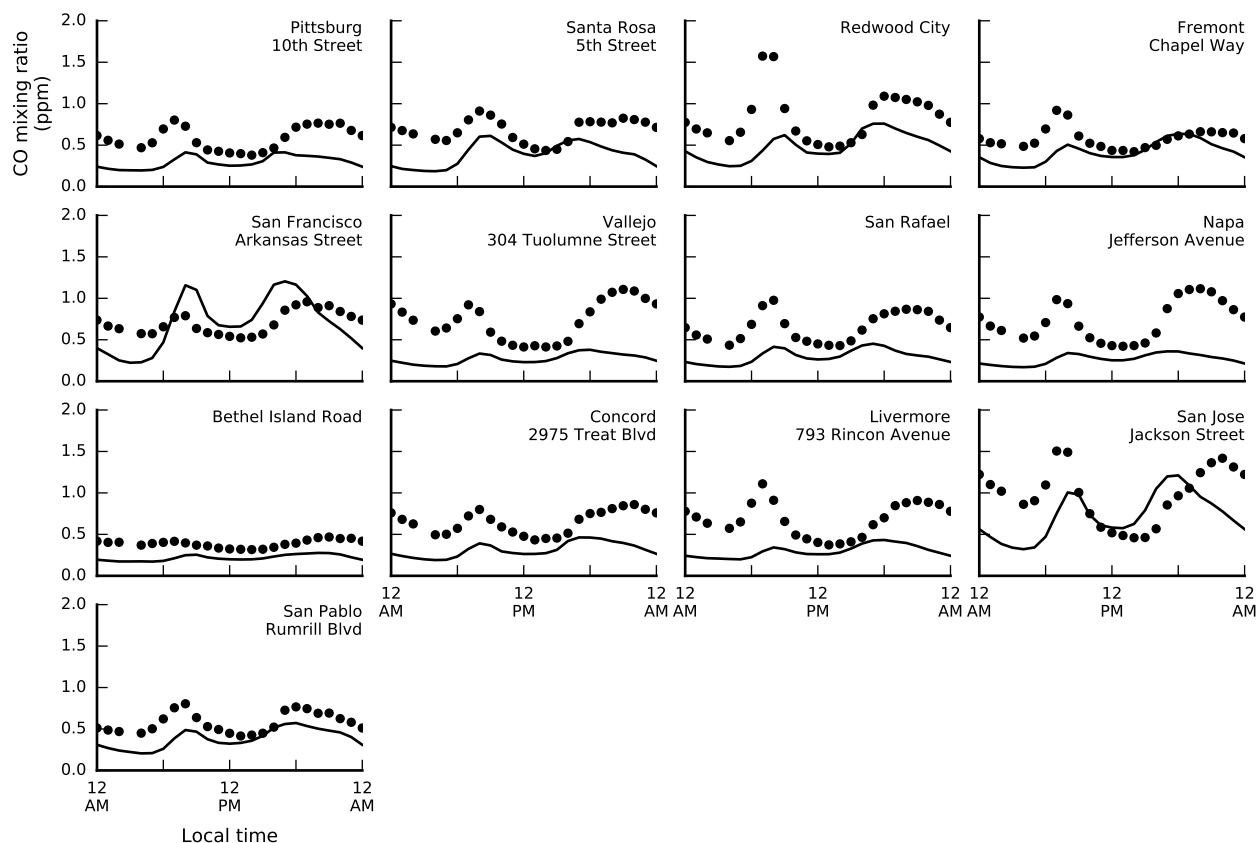


Figure S13: Observed (filled circles) and modeled (solid line) average hourly CO mixing ratios at thirteen different monitoring sites in December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. At each hour, the value is averaged over all the days included in the data set. The locations of the measuring sites are shown on Figure S1.

Table S1: Summary of the comparison between modeled and observed average daily profiles. In each case, the table indicates the average, minimum, and maximum of the mean bias and root mean square error (RMSE) calculated for each of the locations where observations are available. All values are in ppb except for CO (ppm).

	July		December	
	Mean bias	RMSE	Mean bias	RMSE
O ₃	3.8 [−4.4 ; 10.5]	6.4 [3.5 ; 12.7]	6.1 [−6.9 ; 19.6]	9.4 [2.8 ; 19.8]
O ₃ + NO ₂	3.1 [−2.3 ; 7.0]	4.7 [2.3 ; 7.6]	8.9 [4.3 ; 15.4]	9.2 [4.9 ; 15.6]
NO ₂	−0.5 [−5.7 ; 8.8]	4.6 [2.7 ; 9.5]	2.4 [−4.2 ; 13.5]	5.7 [2.8 ; 15.1]
NO	0.7 [−2.8 ; 15.0]	4.6 [1.1 ; 27.5]	−11.8 [−29.7 ; 38.9]	25.8 [8.5 ; 67.6]
NO + NO ₂	0.2 [−8.4 ; 23.8]	8.6 [3.7 ; 35.5]	−9.5 [−31.8 ; 52.3]	27.7 [8.8 ; 80.9]
CO	−0.14 [−0.25 ; −0.02]	0.15 [0.07 ; 0.25]	−0.28 [−0.46 ; 0.01]	0.35 [0.18 ; 0.51]

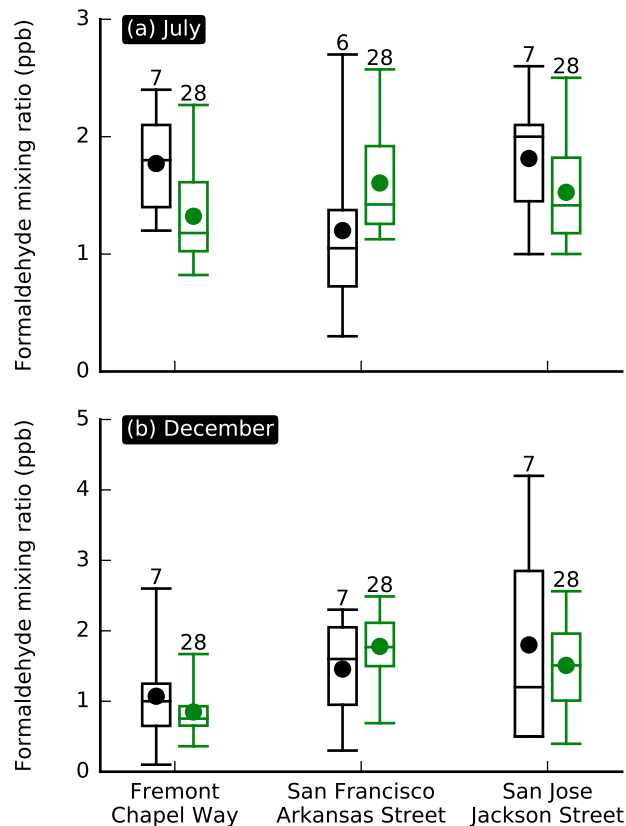


Figure S14: Observed (black, left) and modeled (green, right) 24-hour average formaldehyde mixing ratios at three different monitoring sites in (a) July and (b) December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. Each box-and-whiskers plot represents, for the corresponding data set, the mean (dot), the median (central horizontal line), the first and third quartiles (bottom and top edges of the box), the minimum and maximum values (bottom and top whiskers), and the number of available data points (top number). The locations of the measuring sites are shown on Figure S1.

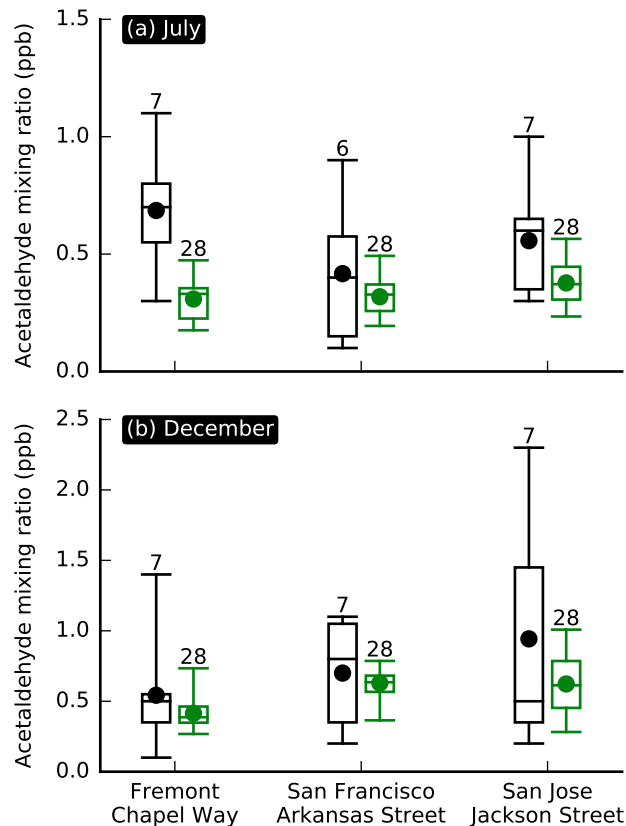


Figure S15: Observed (black, left) and modeled (green, right) 24-hour average acetaldehyde mixing ratios at three different monitoring sites in (a) July and (b) December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. Each box-and-whiskers plot represents, for the corresponding data set, the mean (dot), the median (central horizontal line), the first and third quartiles (bottom and top edges of the box), the minimum and maximum values (bottom and top whiskers), and the number of available data points (top number). The locations of the measuring sites are shown on Figure S1.

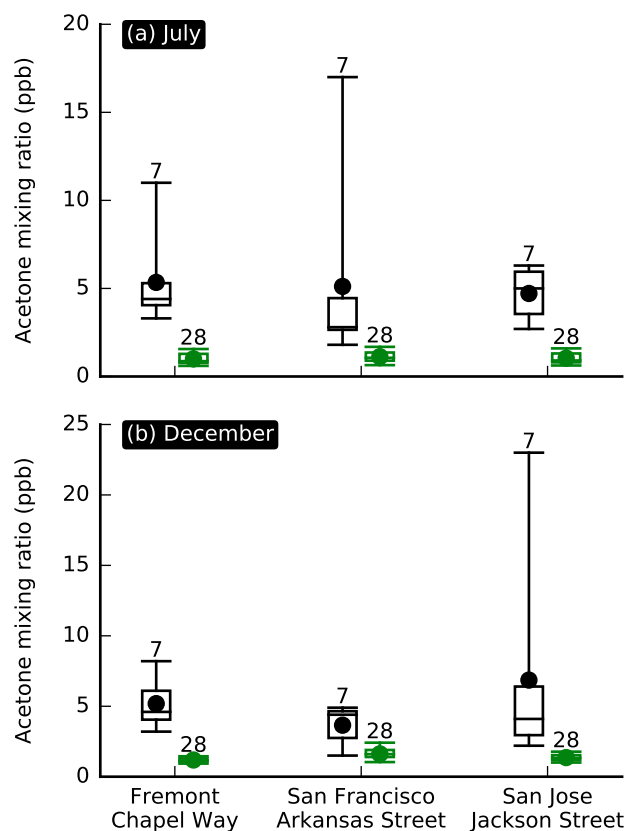


Figure S16: Observed (black, left) and modeled (green, right) 24-hour average acetone mixing ratios at three different monitoring sites in (a) July and (b) December. Observed values combine data from years 2004, 2005, and 2006, by selecting, in each of these years, the days of the year that correspond to the month-long simulation period. Each box-and-whiskers plot represents, for the corresponding data set, the mean (dot), the median (central horizontal line), the first and third quartiles (bottom and top edges of the box), the minimum and maximum values (bottom and top whiskers), and the number of available data points (top number). The locations of the measuring sites are shown on Figure S1.

Table S2: Average surface mixing ratios in the air basin-wide receptor area for each model species, separately for the July and December simulation periods. Surface mixing ratios are mixing ratios in the bottom model layer. Mixing ratios are averaged over the duration of the corresponding simulation period and over the extent of the receptor area. See Carter (2000, Appendix A, Table A-1) for a description of model species.

Model Species	July Simulation	December Simulation	Model Species	July Simulation	December Simulation
NO2	3.8 ppb	10 ppb	BZCO_O2	7.1×10^{-4} ppt	4.1×10^{-5} ppt
NO	1.3 ppb	4.0 ppb	PBZN	0.73 ppt	0.46 ppt
O3P	2.7×10^{-4} ppt	1.0×10^{-4} ppt	BZ_O	4.9×10^{-4} ppt	4.3×10^{-4} ppt
O3	31 ppb	27 ppb	MA_RCO3	0.037 ppt	6.7×10^{-4} ppt
NO3	0.66 ppt	2.2 ppt	MA_PAN	21 ppt	5.5 ppt
N2O5	6.0 ppt	55 ppt	TBU_O	9.9×10^{-7} ppt	5.2×10^{-7} ppt
HNO3	510 ppt	460 ppt	ACET	1000 ppt	1.1 ppb
O1D2	6.1×10^{-10} ppt	4.1×10^{-11} ppt	NPHE	5.1 ppt	7.3 ppt
HO	0.13 ppt	0.010 ppt	PHEN	0.83 ppt	26 ppt
HONO	14 ppt	5.6 ppt	BZNO2_O	2.3×10^{-18} ppt	2.9×10^{-17} ppt
HO2	5.7 ppt	0.82 ppt	HOCOO	1.2×10^{-4} ppt	1.6×10^{-5} ppt
CO	130 ppb	240 ppb	HCOOH	14 ppt	12 ppt
HNO4	4.6 ppt	10 ppt	RCHO	160 ppt	170 ppt
HO2H	1.7 ppb	230 ppt	GLY	11 ppt	15 ppt
SO2	370 ppt	370 ppt	MGLY	35 ppt	25 ppt
SULF	340 ppt	290 ppt	BACL	2.1 ppt	3.5 ppt
C_O2	1.8 ppt	0.43 ppt	CRES	4.8 ppt	5.9 ppt
HCHO	1.3 ppb	550 ppt	BALD	4.3 ppt	8.3 ppt
COOH	53 ppt	4.8 ppt	METHACRO	150 ppt	37 ppt
MEOH	8.1 ppb	510 ppt	MVK	220 ppt	44 ppt
RO2_R	4.2 ppt	2.8 ppt	ISOPROD	46 ppt	26 ppt
ROOH	100 ppt	38 ppt	DCB1	13 ppt	30 ppt
R2O2	5.2 ppt	6.1 ppt	DCB2	0.92 ppt	1.7 ppt
RO2_N	1.6 ppt	1.9 ppt	DCB3	0.38 ppt	0.94 ppt
RNO3	78 ppt	82 ppt	ETHENE	240 ppt	750 ppt
MEK	210 ppt	470 ppt	ISOPRENE	220 ppt	160 ppt
PROD2	150 ppt	430 ppt	TERP	75 ppt	83 ppt
CCO_O2	0.30 ppt	0.020 ppt	ALK1	1.1 ppb	2.3 ppb
PAN	180 ppt	170 ppt	ALK2	360 ppt	910 ppt
CCO_OOH	24 ppt	0.95 ppt	ALK3	620 ppt	1.8 ppb
CCO_OH	25 ppt	9.7 ppt	ALK4	550 ppt	1.7 ppb
RCO_O2	0.079 ppt	2.9×10^{-3} ppt	ALK5	730 ppt	1.9 ppb
PAN2	44 ppt	18 ppt	ARO1	250 ppt	710 ppt
CCHO	220 ppt	320 ppt	ARO2	170 ppt	810 ppt
RCO_OOH	10 ppt	0.19 ppt	OLE1	88 ppt	270 ppt
RCO_OH	21 ppt	21 ppt	OLE2	77 ppt	220 ppt

Table S3: Average surface mixing ratios in the Pittsburgh/Antioch receptor area for each model species, separately for the July and December simulation periods. Surface mixing ratios are mixing ratios in the bottom model layer. Mixing ratios are averaged over the duration of the corresponding simulation period and over the extent of the receptor area. See Carter (2000, Appendix A, Table A-1) for a description of model species.

Model Species	July Simulation	December Simulation	Model Species	July Simulation	December Simulation
NO2	5.1 ppb	11 ppb	BZCO_O2	8.4×10^{-04} ppt	2.6×10^{-05} ppt
NO	1.0 ppb	1.9 ppb	PBZN	1.2 ppt	0.50 ppt
O3P	3.0×10^{-04} ppt	8.4×10^{-05} ppt	BZ_O	4.6×10^{-04} ppt	6.5×10^{-04} ppt
O3	32 ppb	26 ppb	MA_RCO3	0.014 ppt	2.0×10^{-04} ppt
NO3	0.71 ppt	1.1 ppt	MA_PAN	25 ppt	5.0 ppt
N2O5	5.6 ppt	46 ppt	TBU_O	1.2×10^{-06} ppt	6.1×10^{-07} ppt
HNO3	660 ppt	510 ppt	ACET	1.0 ppb	1.1 ppb
O1D2	6.8×10^{-10} ppt	3.8×10^{-11} ppt	NPHE	8.3 ppt	13 ppt
HO	0.17 ppt	9.1×10^{-03} ppt	PHEN	1.5 ppt	68 ppt
HONO	18 ppt	3.8 ppt	BZNO2_O	2.5×10^{-18} ppt	3.7×10^{-17} ppt
HO2	4.7 ppt	0.42 ppt	HOCOO	9.5×10^{-05} ppt	1.0×10^{-05} ppt
CO	140 ppb	250 ppb	HCOOH	19 ppt	10 ppt
HNO4	5.5 ppt	9.0 ppt	RCHO	210 ppt	180 ppt
HO2H	1.7 ppb	220 ppt	GLY	13 ppt	13 ppt
SO2	360 ppt	350 ppt	MGLY	40 ppt	25 ppt
SULF	320 ppt	310 ppt	BACL	2.7 ppt	3.5 ppt
C_O2	1.2 ppt	0.099 ppt	CRES	5.4 ppt	5.9 ppt
HCHO	1.5 ppb	520 ppt	BALD	5.7 ppt	8.4 ppt
COOH	61 ppt	5.0 ppt	METHACRO	140 ppt	32 ppt
MEOH	8.3 ppb	520 ppt	MVK	200 ppt	38 ppt
RO2_R	1.5 ppt	0.21 ppt	ISOPROD	57 ppt	18 ppt
ROOH	96 ppt	29 ppt	DCB1	15 ppt	31 ppt
R2O2	0.98 ppt	0.18 ppt	DCB2	1.2 ppt	1.8 ppt
RO2_N	0.31 ppt	0.045 ppt	DCB3	0.54 ppt	1.1 ppt
RNO3	110 ppt	63 ppt	ETHENE	280 ppt	670 ppt
MEK	240 ppt	470 ppt	ISOPRENE	140 ppt	36 ppt
PROD2	150 ppt	420 ppt	TERP	83 ppt	20 ppt
CCO_O2	0.20 ppt	8.6×10^{-03} ppt	ALK1	1.2 ppb	2.4 ppb
PAN	200 ppt	180 ppt	ALK2	590 ppt	1.2 ppb
CCO_OOH	32 ppt	1.0 ppt	ALK3	810 ppt	1.8 ppb
CCO_OH	33 ppt	9.2 ppt	ALK4	830 ppt	2.0 ppb
RCO_O2	0.074 ppt	1.2×10^{-03} ppt	ALK5	970 ppt	2.0 ppb
PAN2	67 ppt	17 ppt	ARO1	300 ppt	690 ppt
CCHO	280 ppt	310 ppt	ARO2	210 ppt	1.2 ppb
RCO_OOH	14 ppt	0.19 ppt	OLE1	110 ppt	240 ppt
RCO_OH	27 ppt	18 ppt	OLE2	86 ppt	190 ppt

Table S4: Average surface mixing ratios in the San Jose receptor area for each model species, separately for the July and December simulation periods. Surface mixing ratios are mixing ratios in the bottom model layer. Mixing ratios are averaged over the duration of the corresponding simulation period and over the extent of the receptor area. See Carter (2000, Appendix A, Table A-1) for a description of model species.

Model Species	July Simulation	December Simulation	Model Species	July Simulation	December Simulation
NO2	14 ppb	32 ppb	BZCO_O2	1.6×10^{-04} ppt	3.7×10^{-06} ppt
NO	5.4 ppb	31 ppb	PBZN	1.1 ppt	0.61 ppt
O3P	6.0×10^{-04} ppt	2.7×10^{-04} ppt	BZ_O	1.4×10^{-04} ppt	3.6×10^{-05} ppt
O3	24 ppb	7.7 ppb	MA_RCO3	2.2×10^{-03} ppt	4.9×10^{-05} ppt
NO3	0.32 ppt	0.15 ppt	MA_PAN	20 ppt	6.2 ppt
N2O5	7.5 ppt	12 ppt	TBU_O	1.5×10^{-06} ppt	5.4×10^{-07} ppt
HNO3	880 ppt	670 ppt	ACET	1.0 ppb	1.3 ppb
O1D2	5.4×10^{-10} ppt	1.9×10^{-11} ppt	NPHE	8.7 ppt	11 ppt
HO	0.075 ppt	3.4×10^{-03} ppt	PHEN	1.7 ppt	86 ppt
HONO	34 ppt	23 ppt	BZNO2_O	7.9×10^{-19} ppt	1.2×10^{-18} ppt
HO2	0.82 ppt	0.068 ppt	HOCOO	2.0×10^{-05} ppt	1.7×10^{-06} ppt
CO	230 ppb	610 ppb	HCOOH	19 ppt	22 ppt
HNO4	4.3 ppt	3.1 ppt	RCHO	260 ppt	380 ppt
HO2H	1.6 ppb	210 ppt	GLY	20 ppt	31 ppt
SO2	490 ppt	740 ppt	MGLY	45 ppt	42 ppt
SULF	360 ppt	270 ppt	BACL	3.5 ppt	5.2 ppt
C_O2	0.17 ppt	0.014 ppt	CRES	11 ppt	15 ppt
HCHO	1.5 ppb	1.3 ppb	BALD	9.2 ppt	16 ppt
COOH	47 ppt	5.9 ppt	METHACRO	110 ppt	39 ppt
MEOH	7.3 ppb	720 ppt	MVK	170 ppt	47 ppt
RO2_R	0.33 ppt	0.042 ppt	ISOPROD	49 ppt	40 ppt
ROOH	76 ppt	49 ppt	DCB1	26 ppt	52 ppt
R2O2	0.19 ppt	0.050 ppt	DCB2	1.9 ppt	2.7 ppt
RO2_N	0.061 ppt	0.015 ppt	DCB3	0.84 ppt	1.6 ppt
RNO3	130 ppt	140 ppt	ETHENE	790 ppt	2.7 ppb
MEK	250 ppt	530 ppt	ISOPRENE	190 ppt	390 ppt
PROD2	180 ppt	440 ppt	TERP	87 ppt	150 ppt
CCO_O2	0.033 ppt	1.6×10^{-03} ppt	ALK1	1.1 ppb	3.2 ppb
PAN	190 ppt	150 ppt	ALK2	670 ppt	2.1 ppb
CCO_OOH	23 ppt	1.1 ppt	ALK3	1.9 ppb	6.1 ppb
CCO_OH	28 ppt	19 ppt	ALK4	1.7 ppb	5.7 ppb
RCO_O2	0.011 ppt	2.1×10^{-04} ppt	ALK5	2.2 ppb	6.4 ppb
PAN2	54 ppt	21 ppt	ARO1	810 ppt	2.6 ppb
CCHO	370 ppt	570 ppt	ARO2	550 ppt	2.5 ppb
RCO_OOH	10 ppt	0.24 ppt	OLE1	360 ppt	1.2 ppb
RCO_OH	30 ppt	41 ppt	OLE2	310 ppt	1.1 ppb

Table S5: Average surface mixing ratios in the East Bay receptor area for each model species, separately for the July and December simulation periods. Surface mixing ratios are mixing ratios in the bottom model layer. Mixing ratios are averaged over the duration of the corresponding simulation period and over the extent of the receptor area. See Carter (2000, Appendix A, Table A-1) for a description of model species.

Model Species	July Simulation	December Simulation	Model Species	July Simulation	December Simulation
NO2	14 ppb	27 ppb	BZCO_O2	6.4×10^{-05} ppt	3.2×10^{-06} ppt
NO	8.2 ppb	23 ppb	PBZN	0.68 ppt	0.50 ppt
O3P	6.6×10^{-04} ppt	2.1×10^{-04} ppt	BZ_O	8.9×10^{-05} ppt	5.1×10^{-05} ppt
O3	21 ppb	12 ppb	MA_RCO3	1.1×10^{-03} ppt	4.8×10^{-05} ppt
NO3	0.32 ppt	0.27 ppt	MA_PAN	18 ppt	5.3 ppt
N2O5	6.4 ppt	21 ppt	TBU_O	1.3×10^{-06} ppt	4.4×10^{-07} ppt
HNO3	630 ppt	490 ppt	ACET	990 ppt	1.1 ppb
O1D2	4.2×10^{-10} ppt	2.1×10^{-11} ppt	NPHE	5.2 ppt	9.9 ppt
HO	0.055 ppt	3.9×10^{-03} ppt	PHEN	1.6 ppt	62 ppt
HONO	37 ppt	13 ppt	BZNO2_O	6.2×10^{-19} ppt	2.7×10^{-18} ppt
HO2	0.46 ppt	0.099 ppt	HOCOO	1.3×10^{-05} ppt	2.7×10^{-06} ppt
CO	210 ppb	420 ppb	HCOOH	14 ppt	15 ppt
HNO4	3.5 ppt	4.8 ppt	RCHO	210 ppt	260 ppt
HO2H	1.6 ppb	210 ppt	GLY	16 ppt	21 ppt
SO2	1.0 ppb	1.1 ppb	MGLY	39 ppt	31 ppt
SULF	350 ppt	280 ppt	BACL	3.1 ppt	4.0 ppt
C_O2	0.097 ppt	0.021 ppt	CRES	10 ppt	9.4 ppt
HCHO	1.3 ppb	980 ppt	BALD	7.6 ppt	11 ppt
COOH	35 ppt	4.9 ppt	METHACRO	130 ppt	36 ppt
MEOH	7.2 ppb	590 ppt	MVK	190 ppt	42 ppt
RO2_R	0.23 ppt	0.068 ppt	ISOPROD	40 ppt	25 ppt
ROOH	68 ppt	34 ppt	DCB1	24 ppt	36 ppt
R2O2	0.21 ppt	0.097 ppt	DCB2	1.6 ppt	2.1 ppt
RO2_N	0.064 ppt	0.028 ppt	DCB3	0.66 ppt	1.2 ppt
RNO3	89 ppt	93 ppt	ETHENE	860 ppt	2.0 ppb
MEK	210 ppt	480 ppt	ISOPRENE	130 ppt	240 ppt
PROD2	160 ppt	420 ppt	TERP	68 ppt	120 ppt
CCO_O2	0.015 ppt	1.8×10^{-03} ppt	ALK1	1.1 ppb	2.6 ppb
PAN	170 ppt	170 ppt	ALK2	570 ppt	1.5 ppb
CCO_OOH	16 ppt	0.97 ppt	ALK3	1.6 ppb	3.8 ppb
CCO_OH	20 ppt	13 ppt	ALK4	1.4 ppb	3.5 ppb
RCO_O2	4.2×10^{-03} ppt	2.3×10^{-04} ppt	ALK5	2.0 ppb	4.2 ppb
PAN2	36 ppt	18 ppt	ARO1	690 ppt	1.5 ppb
CCHO	300 ppt	440 ppt	ARO2	500 ppt	1.7 ppb
RCO_OOH	6.5 ppt	0.18 ppt	OLE1	290 ppt	710 ppt
RCO_OH	21 ppt	27 ppt	OLE2	270 ppt	660 ppt

Table S6: Overall semi-normalized sensitivities (ppt) of formaldehyde model responses to emissions, by model species. Sensitivities are summed over the entire modeling domain and duration of the simulation period. Species are omitted if the corresponding sensitivities to emissions are negligible for all model responses. For each sensitivity S^\times , the number in parentheses is $S^\times/|S^\times_{\text{tot}}|$ where S^\times_{tot} is the overall semi-normalized emissions sensitivity of the corresponding response, summed across all model species. See Carter (2000, Appendix A, Table A-1) for a description of model species.

Model species	Sensitivity of the Air Basin response		Sensitivity of the Pittsburgh/Antioch response		Sensitivity of the San Jose response		Sensitivity of the East Bay response	
	July	December	July	December	July	December	July	December
NO ₂	-3.3 (-1.1%)	-2.4 (-0.39%)	-2.4 (-0.75%)	-0.93 (-0.32%)	-6.4 (-1.4%)	-4.6 (-0.44%)	-5.0 (-1.3%)	-1.9 (-0.31%)
NO	-55 (-18%)	-77 (-13%)	-36 (-12%)	-28 (-9.7%)	-120 (-26%)	-150 (-14%)	-87 (-23%)	-60 (-10%)
CO	-0.91 (-0.30%)	0.0 (0.0%)	-1.7 (-0.54%)	0.0 (0.0%)	-0.78 (-0.17%)	0.0 (0.0%)	-0.84 (-0.22%)	0.0 (0.0%)
HCHO	170 (54%)	540 (88%)	86 (27%)	250 (85%)	270 (58%)	920 (90%)	280 (75%)	540 (90%)
MEOH	0.0 (0.0%)	0.0 (0.0%)	0.39 (0.12%)	0.0 (0.0%)	0.47 (0.10%)	0.0 (0.0%)	0.0 (0.0%)	0.0 (0.0%)
CCHO	1.7 (0.56%)	0.92 (0.15%)	1.9 (0.61%)	0.0 (0.0%)	2.8 (0.62%)	2.2 (0.21%)	2.5 (0.65%)	0.64 (0.11%)
ACET	0.0 (0.0%)	0.0 (0.0%)	0.34 (0.11%)	0.0 (0.0%)	0.51 (0.11%)	0.0 (0.0%)	0.0 (0.0%)	0.0 (0.0%)
RCHO	0.0 (0.0%)	0.0 (0.0%)	0.39 (0.12%)	0.0 (0.0%)	0.54 (0.12%)	0.0 (0.0%)	0.0 (0.0%)	0.0 (0.0%)
ETHENE	34 (11%)	20 (3.4%)	41 (13%)	10 (3.6%)	54 (12%)	36 (3.5%)	44 (12%)	16 (2.7%)
ISOPRENE	72 (23%)	29 (4.8%)	120 (37%)	10 (3.6%)	100 (22%)	45 (4.4%)	44 (11%)	23 (3.8%)
TERP	14 (4.7%)	12 (1.9%)	17 (5.3%)	3.8 (1.3%)	21 (4.6%)	19 (1.8%)	12 (3.1%)	9.6 (1.6%)
ALK1	0.50 (0.16%)	0.0 (0.0%)	1.0 (0.32%)	0.0 (0.0%)	0.99 (0.22%)	0.0 (0.0%)	0.0 (0.0%)	0.0 (0.0%)
ALK2	0.61 (0.20%)	0.0 (0.0%)	0.80 (0.25%)	0.0 (0.0%)	1.5 (0.33%)	0.0 (0.0%)	0.0 (0.0%)	0.0 (0.0%)
ALK3	5.2 (1.7%)	1.2 (0.20%)	8.3 (2.6%)	0.59 (0.20%)	9.3 (2.0%)	2.3 (0.22%)	4.4 (1.2%)	0.97 (0.16%)
ALK4	3.9 (1.3%)	0.0 (0.0%)	7.4 (2.4%)	0.0 (0.0%)	7.7 (1.7%)	0.0 (0.0%)	2.4 (0.63%)	0.0 (0.0%)
ALK5	3.3 (1.1%)	-1.7 (-0.28%)	7.9 (2.5%)	-0.69 (-0.24%)	8.7 (1.9%)	-3.1 (-0.30%)	0.0 (0.0%)	-1.3 (-0.21%)
ARO1	3.1 (1.0%)	1.0 (0.17%)	4.3 (1.4%)	0.0 (0.0%)	6.2 (1.4%)	2.3 (0.23%)	2.7 (0.72%)	0.65 (0.11%)
ARO2	9.9 (3.2%)	8.5 (1.4%)	12 (3.8%)	3.1 (1.1%)	17 (3.8%)	18 (1.7%)	11 (3.0%)	6.1 (1.0%)
OLE1	18 (5.9%)	17 (2.8%)	19 (6.1%)	8.7 (3.0%)	31 (6.7%)	30 (2.9%)	22 (5.9%)	13 (2.2%)
OLE2	33 (11%)	62 (10%)	32 (10%)	34 (12%)	56 (12%)	110 (10%)	40 (11%)	50 (8.3%)
Total	310	610	310	290	460	1000	380	600
Categories:								
NO _x	-59 (-19%)	-79 (-13%)	-39 (-12%)	-29 (-10%)	-130 (-28%)	-150 (-15%)	-92 (-24%)	-62 (-10%)
VOCs (all)	370 (120%)	690 (110%)	360 (110%)	320 (110%)	590 (130%)	1200 (110%)	470 (120%)	660 (110%)
VOCs (antropogenic)	280 (90%)	640 (110%)	220 (70%)	300 (100%)	460 (100%)	1100 (110%)	410 (110%)	630 (100%)
VOCs (biogenic)	89 (29%)	43 (7.1%)	140 (43%)	16 (5.4%)	130 (28%)	67 (6.5%)	57 (15%)	34 (5.7%)
VOCs (on-road)	110 (35%)	220 (35%)	92 (29%)	86 (30%)	190 (40%)	400 (39%)	150 (41%)	230 (39%)

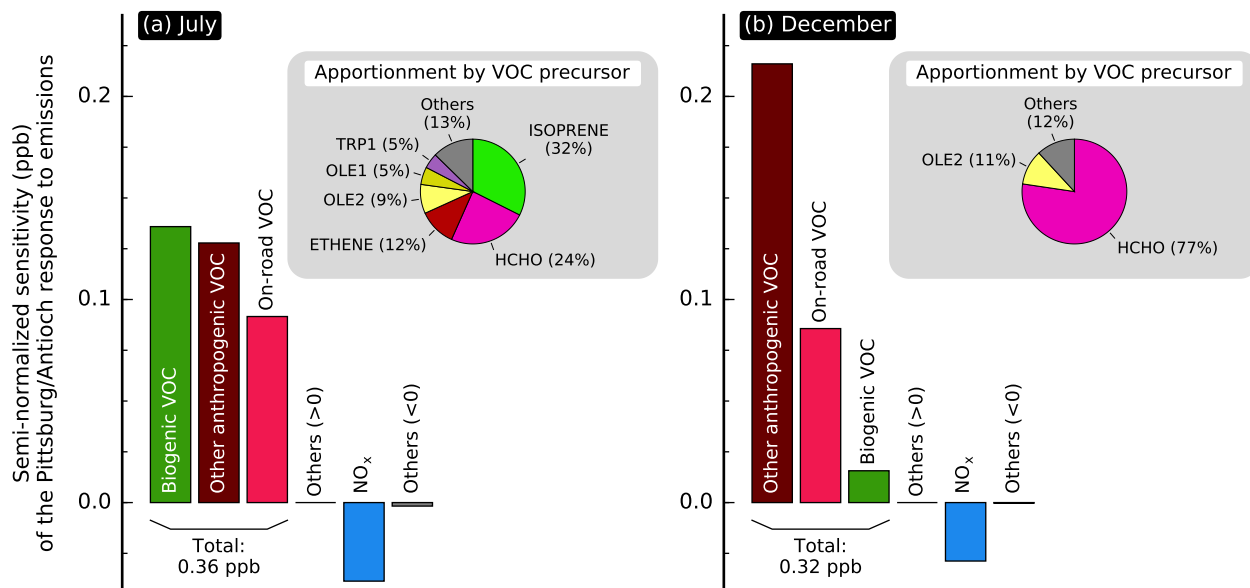


Figure S17: Semi-normalized sensitivity of the Pittsburgh/Antioch population-weighted formaldehyde mixing ratio to emissions, in (a) July and (b) December. An apportionment by precursor species of the sensitivity to overall VOC emissions is shown in the pie chart insets.

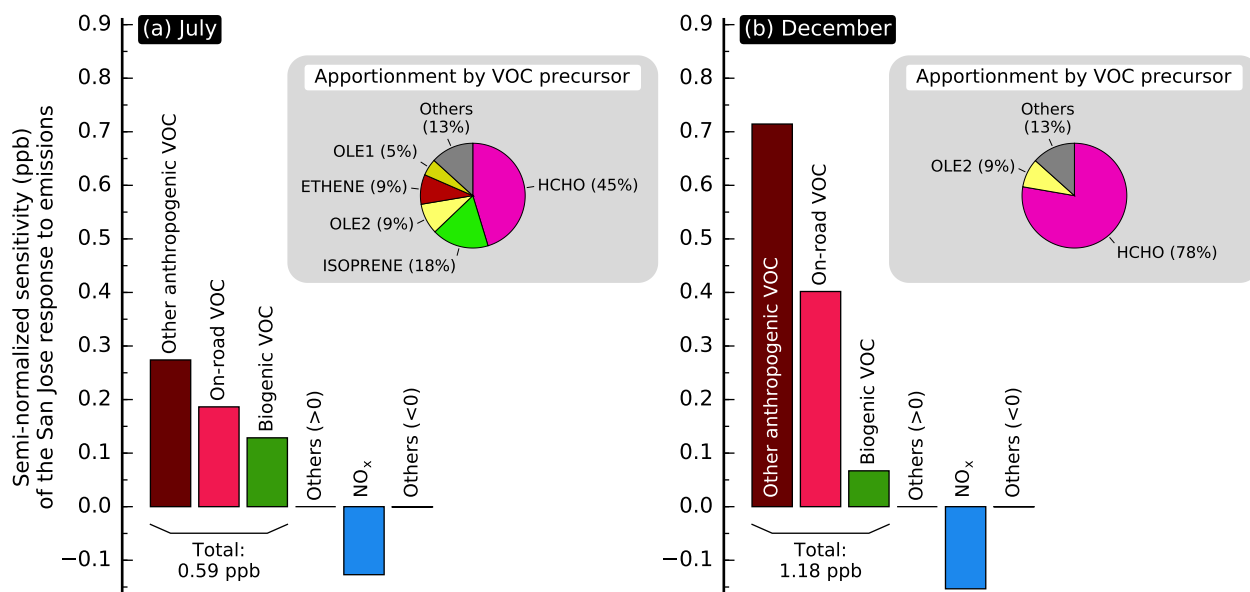


Figure S18: Semi-normalized sensitivity of the San Jose population-weighted formaldehyde mixing ratio to emissions, in (a) July and (b) December. An apportionment by precursor species of the sensitivity to overall VOC emissions is shown in the pie chart insets.

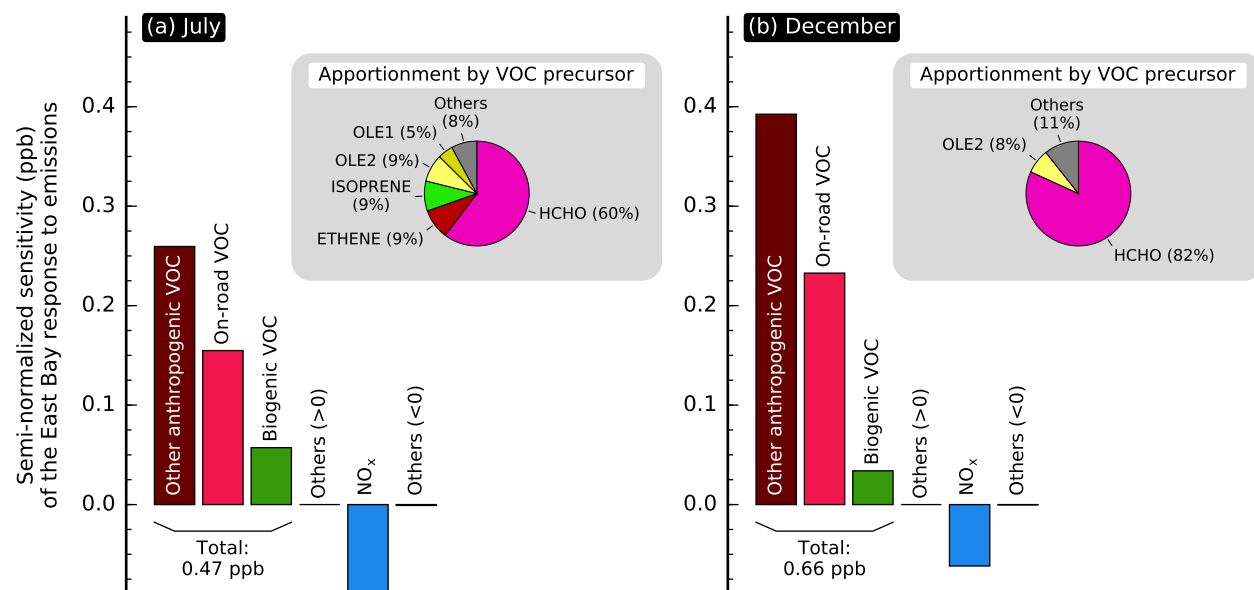


Figure S19: Semi-normalized sensitivity of the East Bay population-weighted formaldehyde mixing ratio to emissions, in (a) July and (b) December. An apportionment by precursor species of the sensitivity to overall VOC emissions is shown in the pie chart insets.

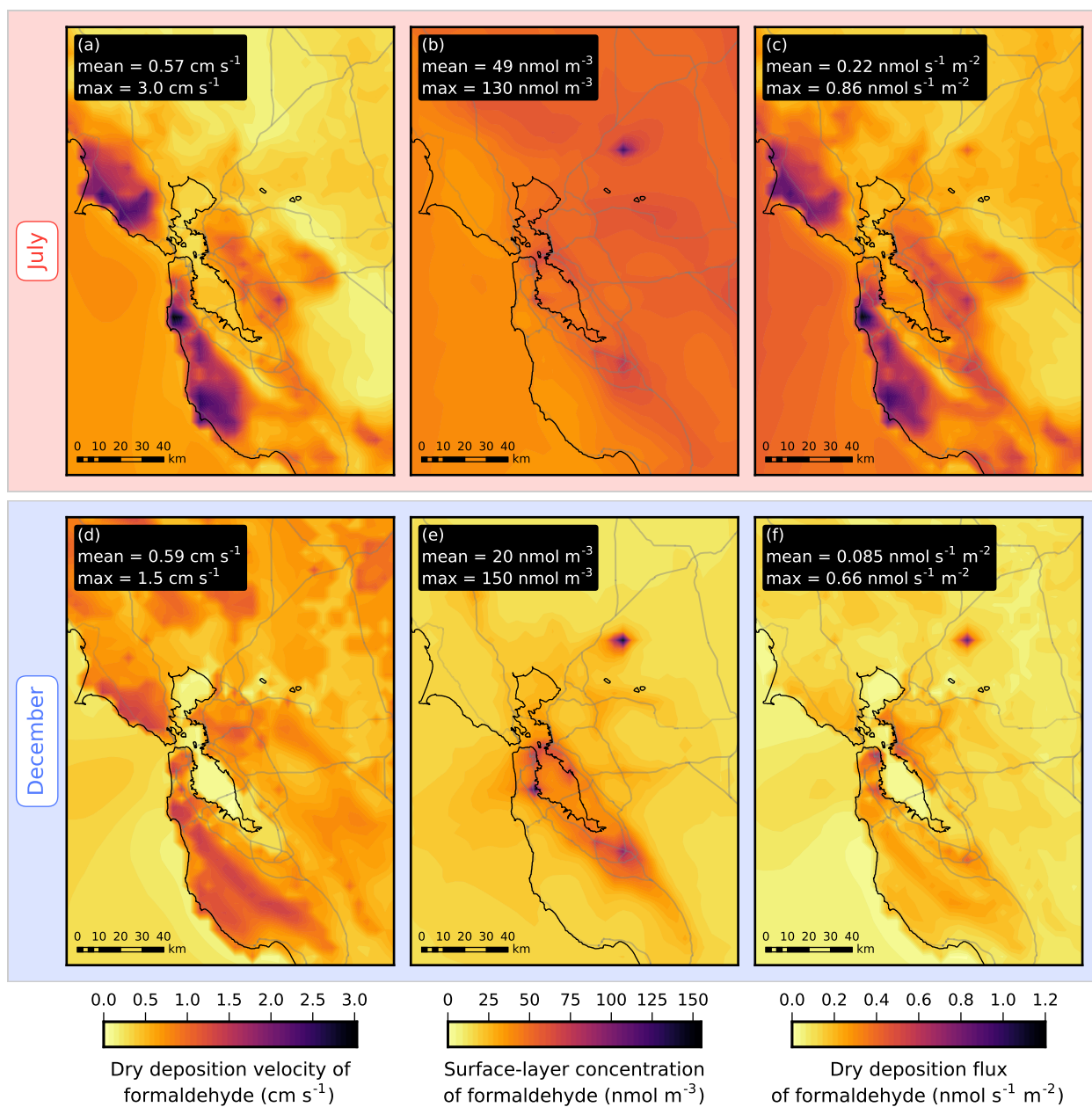


Figure S20: Dry deposition velocity (v_d), surface-layer concentration (C), and dry deposition flux (F) of formaldehyde, where $F = v_d \times C$. Dry deposition velocities were calculated by the MM5 model (BAAQMD, 2009).

References

- BAAQMD (Bay Area Air Quality Management District): Toxics Modeling to Support the Community Air Risk Evaluation (CARE) Program, Research and Modeling Section Publication No. 200906-002-TX. San Francisco, CA. June 2009, 2009.
- Carter, W. P. L.: Documentation of the SAPRC-99 Chemical Mechanism for VOC reactivity assessment. Volume 1 of 2. Documentation Text, Final Report to California Air Resources Board. Contract 92-329 and Contract 95-308. Report number 00-AP-RT17-001-FR, May 8th 2000, 2000.