



2003 PHOTOCHEMICAL ASSESSMENT MONITORING STATIONS (PAMS)

New Jersey Department of Environmental Protection

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Most ground-level ozone is the result of oxides of nitrogen (NO_x) and volatile organic compounds (VOCs) reacting in the presence of sunlight. As a result, it is necessary to measure these ozone forming pollutants, also known as precursor pollutants, to effectively evaluate strategies for reducing ozone levels. The Photochemical Assessment Monitoring Stations (PAMS) network was established for this purpose. Data from the PAMS network is used to better characterize the nature and extent of the O₃ problem, track VOC and NO_x emission inventory reductions, assess air quality trends, and make attainment/nonattainment decisions. PAMS monitor both criteria and non-criteria pollutants including ozone (O₃), oxides of nitrogen (NO_x), nitric oxide (NO), nitrogen dioxide (NO₂), and specific VOCs, including several carbonyls, that are important in ozone formation. In addition, the measurement of specific weather parameters (e.g. Wind speed/direction, temperature) is required at all PAMS, and upper air weather measurements are required in certain areas. The VOC and carbonyl measurements are only taken during the peak part of the ozone season, from June 1st to August 31st each year.

The PAMS network is designed around metropolitan areas where ozone is a significant problem, and each site in the network has a specific purpose as shown in the Figure 1 below. New Jersey is part of the Philadelphia and New York Metropolitan areas and has a total of three PAMS sites. A Type 3 maximum ozone site for the Philadelphia area is located at Rider University in Mercer County, a Type 2 maximum emissions site is located downwind of the Philadelphia Metropolitan urban area in Camden, and a site at Rutgers University in New Brunswick has been designated both a PAMS Type 1 upwind site for the New York urban area, as well as a Type 4 downwind site for the Philadelphia Metropolitan urban area. An upper air weather monitoring station is also located at the Rutgers University site. All of the PAMS sites for the Philadelphia and New York City areas are shown in Figure 2.

Figure 1

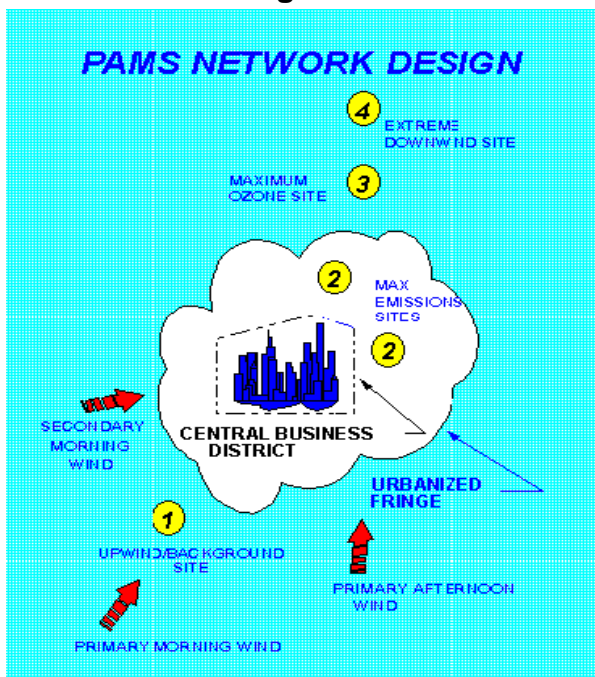


Figure 2

Regional PAMS Sites



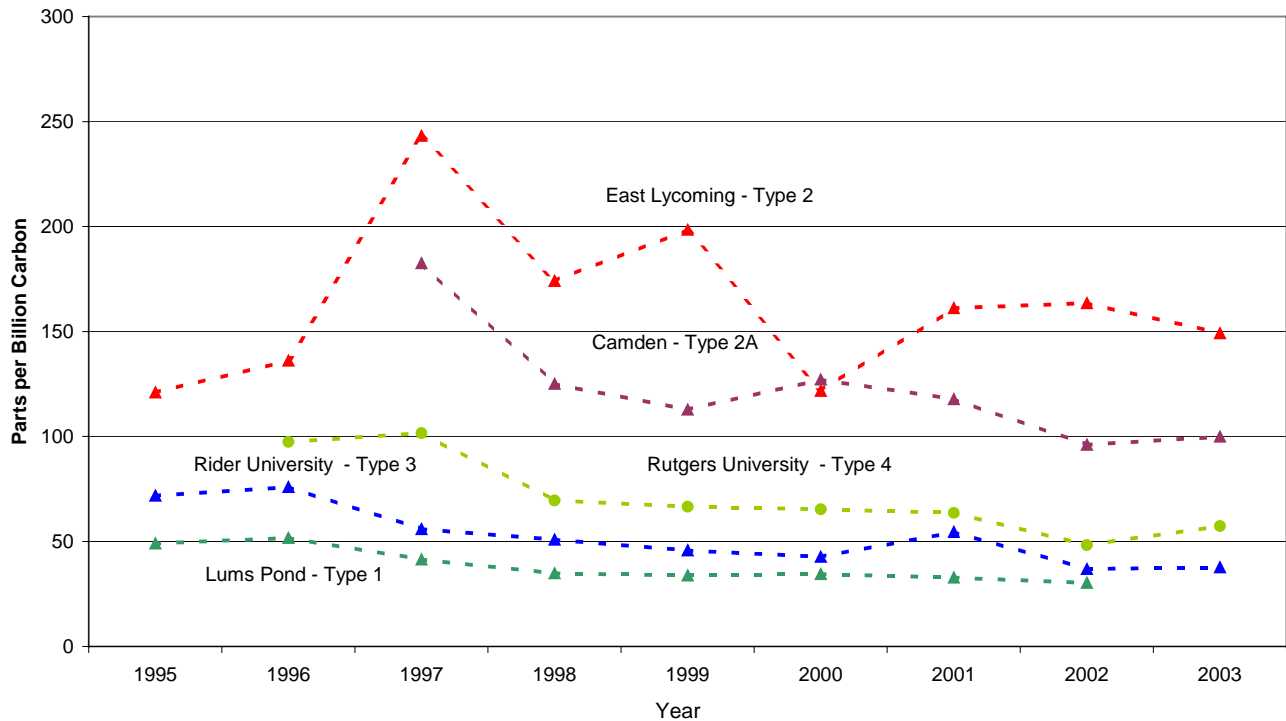
Note: Rutgers University PAMS site is both Type 4 for Philadelphia and Type 1 for New York City.

⁵ USEPA , PAMS General Information

PAMS (CONT.)

Figure 3 shows VOC trends for the PAMS sites in the Philadelphia area. In general, for Lums Pond (upwind - Type 1), Rider University (maximum ozone concentration - Type 3) and Rutgers University (downwind - Type 4), VOCs have declined over the measurement period. The improvements were initially more dramatic, with more level, though still declining concentrations, over the last several years. The maximum emissions -Type 2 sites (Camden and East Lycoming) for this area show more variation from year to year, though the trend at both sites is downward since 1997. Operation of the Lums Pond site was discontinued after the 2002 season by Delaware's Department of Natural Resources and Environmental Control (DNREC).

Figure 3
Philadelphia Region
Total Non-methane Organic Carbon (TNMOC)
Seasonal Average 1995-2003



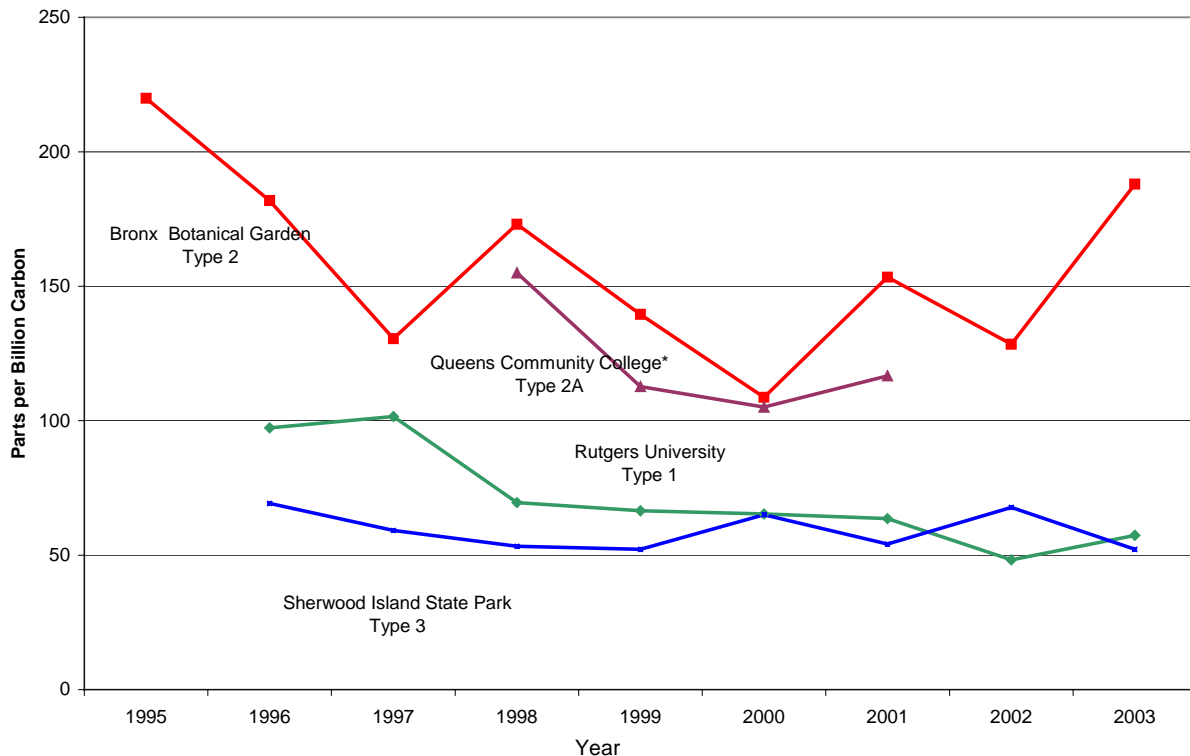
PAMS (cont.)

Figure 4 shows VOC trends for the PAMS sites in the New York City metropolitan area. In general, observations here are similar to those for the Philadelphia area. The Type 2 site in the NY area at the Bronx Botanical Gardens shows even more year to year variability than does the Philadelphia Type 2 site at East Lycoming. Operation of the Queens Community College site was discontinued after the 2001 season.

In conclusion, trends for VOC values measured at all PAMS sites in the Philadelphia and New York City areas show a decline over the time period these measurements were made. Changes in gasoline formulation over the period as well as the effect of newer, cleaner vehicles replacing older vehicles in the automotive fleet could account for the reductions. Type 2 sites, though impacted by vehicle emissions, are also affected by urban stationary sources whose emission trends over the measurement period are less clear, hence these sites seem to show more year to year variability. All sites are also impacted by naturally occurring isoprene, which is emitted by trees. All VOCs are not equal in their contribution to ozone formation and while isoprene levels are generally lower than many other VOCs, isoprene can account for a significant amount of the ozone forming potential, especially at the non-urban sites. Isoprene levels are also highest during the middle of the day, when photochemical conditions are most conducive to ozone formation. Isoprene levels are thought to be influenced by factors that affect tree health and growth, such as rainfall and severe temperatures.

Summaries of results for all the VOCs and carbonyls measured at the New Jersey PAMS sites are provided in Table 1 and Table 2.

Figure 4
New York City Region
Total Non-methane Organic Carbon (TNMOC)
Seasonal Average 1995-2003



* Operations of the Queens Community College site was discontinued after the 2001 season.

Table 1
Summary of Photochemical Assessment Monitoring (PAMS) Data
June, July, and August, 2003

Parts Per Billion (Volume) – ppbv
Parts Per Billion (Carbon) – ppbC
Max – Maximum Avg - Average

	Camden Lab				Rider University				Rutgers University			
	ppbv		ppbC		ppbv		ppbC		ppbv		ppbC	
	Max	Avg	Max	Avg	Max	Avg	Max	Avg	Max	Avg	Max	Avg
Acetylene	5.09	0.39	10.18	0.77	0.62	0.14	1.23	0.28	6.68	0.33	13.36	0.66
Benzene	4.71	0.35	28.25	2.12	0.74	0.15	4.41	0.88	0.77	0.15	4.62	0.88
n-Butane	46.98	1.62	187.93	6.49	22.83	0.42	91.3	1.67	5.90	0.57	23.61	2.27
1-Butene	2.97	0.13	11.86	0.53	0.20	0.03	0.78	0.13	0.77	0.05	3.07	0.19
cis-2-Butene	1.59	0.08	6.35	0.31	0.12	0.02	0.47	0.08	0.54	0.03	2.17	0.12
trans-2-Butene	1.91	0.09	7.62	0.38	0.11	0.02	0.45	0.08	0.54	0.03	2.14	0.12
Cyclohexane	2.49	0.11	14.93	0.65	0.17	0.03	1.00	0.20	0.24	0.04	1.42	0.22
Cyclopentane	1.03	0.10	5.17	0.49	0.23	0.06	1.13	0.28	0.52	0.04	2.61	0.21
n-Decane	4.69	0.06	46.92	0.61	2.13	0.03	21.27	0.27	0.63	0.04	6.31	0.35
m-Diethylbenzene	0.37	0.02	3.7	0.23	0.21	0.02	2.1	0.20	1.23	0.02	12.29	0.15
p-Diethylbenzene	0.49	0.02	4.87	0.25	0.28	0.01	2.81	0.14	0.12	0.02	1.19	0.16
2,2-Dimethylbutane	3.80	0.17	19.01	0.83	0.20	0.04	0.99	0.18	0.97	0.04	4.86	0.22
2,3-Dimethylbutane	1.29	0.15	6.46	0.73	0.25	0.05	1.25	0.27	0.88	0.07	4.39	0.34
2,3-Dimethylpentane	0.50	0.08	3.53	0.55	0.22	0.04	1.51	0.29	0.28	0.04	1.95	0.30
2,4-Dimethylpentane	0.45	0.06	3.17	0.42	0.12	0.03	0.81	0.23	0.21	0.03	1.45	0.24
Ethane	21.54	3.95	43.08	7.90	13.70	2.39	27.39	4.78	43.27	2.87	86.54	5.74
Ethylbenzene	0.66	0.09	5.29	0.70	1.72	0.04	13.72	0.35	0.39	0.05	3.13	0.40
Ethylene (Ethene)	15.24	1.31	30.47	2.63	3.33	0.42	6.66	0.85	19.62	1.16	39.24	2.32
Isomers of Ethyltoluene	4.16	0.13	37.48	1.15	0.77	0.06	6.94	0.53	1.20	0.10	10.72	0.93
n-Heptane	1.43	0.16	10.00	1.11	0.25	0.05	1.77	0.35	0.72	0.06	5.06	0.44
Hexane	4.64	0.34	27.85	2.03	0.63	0.12	3.80	0.69	2.90	0.17	17.37	1.03
1-Hexene	0.75	0.03	4.49	0.18	0.20	0.02	1.19	0.14	0.11	0.02	0.68	0.09
Isobutane	24.74	1.17	98.97	4.67	11.90	0.28	47.59	1.13	3.61	0.36	14.44	1.43
Isopentane	17.71	1.46	88.56	7.30	2.54	0.50	12.70	2.52	13.42	0.68	67.09	3.42
Isoprene	1.43	0.23	7.15	1.17	2.68	0.21	13.42	1.06	4.35	0.59	21.73	2.95
Isopropylbenzene	1.33	0.07	12.01	0.59	0.14	0.02	1.28	0.16	0.08	0.02	0.75	0.14
Methylcyclohexane	2.29	0.12	16.01	0.87	0.27	0.04	1.87	0.27	0.66	0.05	4.61	0.34
Methylcyclopentane	2.39	0.17	14.35	1.03	0.31	0.07	1.84	0.40	0.67	0.08	3.99	0.46
2-Methylheptane	0.43	0.05	3.42	0.39	0.37	0.02	2.93	0.13	0.34	0.02	2.75	0.18
3-Methylheptane	0.46	0.05	3.71	0.36	0.21	0.02	1.71	0.16	0.39	0.02	3.14	0.20
2-Methylhexane	0.86	0.13	6.04	0.90	0.24	0.05	1.71	0.34	0.37	0.06	2.57	0.41

Table 1 (Continued)
Summary of Photochemical Assessment Monitoring (PAMS) Data
June, July, and August, 2003

	Camden Lab				Rider University				Rutgers University			
	ppbv		ppbC		ppbv		ppbC		ppbv		ppbC	
	Max	Avg	Max	Avg	Max	Avg	Max	Avg	Max	Avg	Max	Avg
3-Methylhexane	1.12	0.16	7.83	1.12	0.31	0.06	2.14	0.42	0.42	0.07	2.96	0.49
2-Methylpentane	3.46	0.39	20.76	2.35	0.69	0.14	4.13	0.82	2.25	0.17	13.49	1.03
3-Methylpentane	2.21	0.25	13.27	1.50	0.45	0.09	2.69	0.55	1.38	0.11	8.29	0.67
n-Nonane	2.83	0.06	25.45	0.56	1.11	0.02	10.01	0.22	0.53	0.03	4.79	0.29
n-Octane	0.62	0.08	4.99	0.61	1.04	0.03	8.31	0.20	1.06	0.04	8.47	0.30
n-Pentane	13.89	0.84	69.46	4.20	1.82	0.26	9.11	1.28	4.32	0.37	21.59	1.86
1-Pentene	0.65	0.06	3.23	0.31	0.07	0.02	0.37	0.09	0.36	0.03	1.81	0.15
cis-2-Pentene	0.38	0.04	1.88	0.21	0.06	0.02	0.3	0.08	0.31	0.03	1.55	0.13
trans-2-Pentene	0.63	0.07	3.15	0.34	0.14	0.02	0.69	0.11	0.56	0.04	2.8	0.18
Propane	57.03	3.97	171.1	11.90	20.70	1.35	62.1	4.05	16.70	1.69	50.11	5.08
n-Propylbenzene	0.32	0.03	2.9	0.26	0.16	0.02	1.41	0.14	0.14	0.02	1.23	0.17
Propylene (Propene)	32.20	0.88	96.6	2.64	2.55	0.22	7.64	0.67	4.21	0.34	12.62	1.03
Styrene	0.40	0.04	3.21	0.36	0.30	0.02	2.4	0.18	0.16	0.03	1.28	0.21
Toluene	5.37	0.69	37.59	4.82	33.85	0.35	236.98	2.46	14.13	0.72	98.91	5.04
1,2,3-Trimethylbenzene	1.68	0.05	15.09	0.46	0.73	0.06	6.57	0.52	0.58	0.05	5.23	0.47
1,2,4-Trimethylbenzene	1.79	0.09	16.15	0.79	0.77	0.04	6.96	0.33	0.63	0.06	5.67	0.55
1,3,5-Trimethylbenzene	0.78	0.05	6.99	0.41	0.45	0.03	4.06	0.23	0.27	0.03	2.46	0.24
2,2,4-Trimethylpentane	2.17	0.21	17.35	1.65	0.43	0.08	3.45	0.67	0.82	0.11	6.57	0.92
2,3,4-Trimethylpentane	0.89	0.06	7.09	0.52	0.13	0.03	1.07	0.23	0.28	0.04	2.22	0.31
n-Undecane	2.60	0.04	28.57	0.47	1.79	0.02	19.73	0.22	0.37	0.03	4.07	0.33
m/p-Xylene	2.20	0.24	17.6	1.92	6.29	0.12	50.32	1.00	1.43	0.16	11.42	1.32
o-Xylene	0.78	0.09	6.24	0.75	1.10	0.05	8.8	0.38	0.54	0.06	4.34	0.47

Table 2
Camden Lab
PAMS Carbonyls
June, July, and August, 2003
Parts Per Billion (Volume)
12 Sampling Dates (96 Observations)

	# of Detects*	Max	Avg		# of Detects*	Max	Avg
Acetaldehyde	96	5.70	1.14	Formaldehyde	96	10.02	2.68
Acetone	96	4.87	2.67	Hexaldehyde	52	0.10	0.01
Benzaldehyde	95	0.08	0.03	Isovaleraldehyde	31	0.04	0.01
Butyr/Isobutyraldehyde	61	0.25	0.04	Propionaldehyde	72	0.33	0.05
Crotonaldehyde	94	0.52	0.13	Tolualdehyde	75	0.32	0.04
2,5-Dimethylbenzaldehyde	0	0.00	0.00	Valeraldehyde	93	0.09	0.02

* The number of samples, out of a possible 96, in which the compound was detected.

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