# **Chapter 1 Results of the Initial Environmental Survey in FY 2007**

## 1. Purpose of the survey

The Initial Environmental Survey is aimed at understanding the environmental presence of chemicals requiring examination of the appropriateness of the designation as a Designated Chemical Substance in the Law Concerning Reporting, etc. of Releases of Specific Chemical Substances to the Environment and Promoting Improvement in Their Management (Law No. 86 of 1999) (hereafter, the PRTR Law) and chemicals requiring survey from social viewpoints.

## 2. Target chemicals

In the FY 2007 Initial Environmental Survey, 27chemicals that were selected and designated as target chemicals. The combinations of target chemicals and the surveyed media are given below.

	Target chemicals		Designated Class		Su	rveyed medi	a
No.	Name	The Chemical Substances Control Law	The PRTR Law (Current)	The PRTR Law (New) Effective from October 1, 2009	Surface water	Sediment	Air
[1]	Adipic acid						0
[2]	Ethyleneimine		I	I			0
[3]	4'-Ethoxyacetanilide (synonym:Phenacetin)	II Monitored	I				0
[4]	2,4-Xylenol		II	I	0		
[5]	Quinoline			I	0		
[6]	5-Chloro- <i>N</i> -{2-[4-(2-ethoxyethyl)-2,3-dimethylph enoxy]ethyl}-6-ethylpyrimidine-4-amine (synonym:Pylimidifen)		II				0
[7]	1-Chloronaphthalene	III Monitored	II	II			0
[8]	2-Chloronitrobenzene			I	0		
[9]	Salicylaldehyde		I	I	0		
[10]	2,6-Dinitrotoluene	II Monitored III Monitored	I	I	0		
[11]	<i>m</i> -Dinitrobenzene	II Monitored	II	II	0		
[12]	Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene)	III Monitored	II	II	0	0	
[13]	Dimethyl 4,4'-(o-phenylene)bis(3-thioallophanate) (synonym: Thiophanate-methyl)			I	0		
[14]	Dimethyl terephthalate		I	I			0
[15]	Propylene dinitrate		II				0
[16]	o-Nitroaniline			I	0		
[17]	<i>m</i> -Nitroaniline	II Monitored	II	II	0		
[18]	Vanadium and its compounds (as Vanadium)				0		
[19]	Phenanthrene		II		0	0	
[20]	Phenyloxirane (synonym: Styreneoxide)		I	I	0		
[21]	Dimethyl phthalate				0	0	
[22]	Benzyl alcohol						0
[23]	Methylhydrazine	II Monitored	II	II	0		
[24]	2-Methyl-1,1'-biphenyl-3-ylmethyl (Z)-3-(2-chloro-3,3,3-trifluoro-1-propenyl)-2,2-di methylcyclopropanecarboxylate (synonym: Bifenthrin)	II Monitored	II	II	0		
[25]	2-(1-Methylpropyl)-4,6-dinitrophenol	II Monitored III Monitored	I	I			0
[26]	Mercaptoaceticacid		I		0		
[27]	Triphenyl phosphate			I			0
ONT.	(a) "The Chemical Substances Central Law" hereafte	"T C	· 4 F			C .	,

(Note) "The Chemical Substances Control Law" hereafter means "Law Concerning the Examination and Regulation of Manufacture, etc. Of Chemical Substances (Law No. 117 of 1973)."

Chemical and physical properties of target chemicals of the	Initial Environmental Surv	vey are as follows.
[1] Adipic acid		•
	Molecular formula: CAS:	C <sub>6</sub> H <sub>10</sub> O <sub>4</sub> 124-04-9
0	ENCS:	2-858
	MW:	146.14
НО	mp:	152°C¹)
<u> </u>	bp:	337.5°C¹)
0	SW:	15g/L (15°C) <sup>1)</sup>
	Specific gravity:	1.36 (25/4°C) <sup>1)</sup>
	logPow:	$0.08^{2)}$
[2] Ethyleneimine		
	Molecular formula:	$C_2H_5N$
	CAS:	151-56-4
Н	ENCS:	5-2
Ñ	MW:	43.07
/ \	mp:	$-71.5^{\circ}C^{3}$
	bp:	$55\sim56^{\circ}C^{3)}$
	SW:	Readily soluble <sup>4)</sup>
	Specific gravity: logPow:	$0.83 (24/4^{\circ}C)^{-1}$ - $0.28^{5}$
[3] 4'-Ethoxyacetanilide (synonym: Phenacetin)	logrow:	-U.20
	M 1 1 C 1	C II NO
	Molecular formula: CAS:	C <sub>10</sub> H <sub>13</sub> NO <sub>2</sub> 62-44-2
HNO	ENCS:	3-697
	MW:	179.22
	mp:	134∼135°C¹)
	bp:	242~245°C³)
	SW:	766mg/L (25°C) $^{6)}$
	Specific gravity:	1.36 (20/4°C) <sup>7)</sup>
V	logPow:	$1.58^{8)}$
[4] 2,4-Xylenol		
	Molecular formula:	$C_8C_{10}O$
	CAS:	105-67-9
	ENCS:	3-521, 4-57
	MW: mp:	122.17 24.5℃ <sup>9)</sup>
	bp:	24.5 ℃ (760 mmHg) 9)
	SW:	7.87×10 <sup>3</sup> mg/L $(25^{\circ}C)^{-10}$
<b>Y</b>	Specific gravity:	Uncertain (25 C)
OH	logPow:	2.30 <sup>11)</sup>
[5] Quinoline		
	Molecular formula:	C <sub>9</sub> H <sub>7</sub> N
∧ .N.	CAS:	91-22-5
	ENCS:	5-794
	MW:	129.16
	mp:	-14.78°C <sup>9)</sup>
	bp:	$237.7^{\circ}C^{1)}$
	SW:	$6.11 \times 10^3 \text{ mg/L} (25^{\circ}\text{C})^{-12}$
	Specific gravity:	$1.09 (25/4^{\circ}C)^{-1}$ $2.03^{9}$
Abbreviations) CAS: CAS registry number, ENCS: registry n	logPow:	

(Abbreviations) CAS: CAS registry number, ENCS: registry number in the Existing and New Chemical Substances List, MW: molecular weight, mp: melting point, bp: boiling point, SW: solubility in water, logPow: *n*-octanol-water partition coefficient, kPa: kilopascal (1 atom = 101.3kPa).

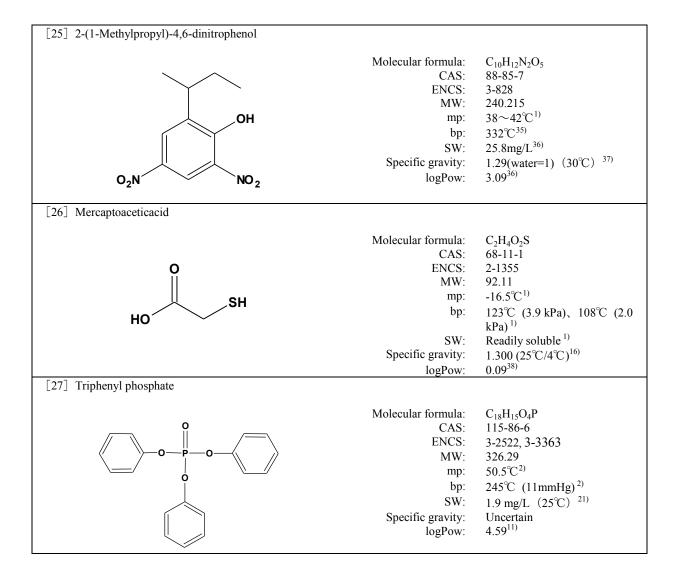
[6] 5-Chloro- <i>N</i> -{2-[4-(2-ethoxyethyl)-2,3-dimethylpheno	xy]ethyl}-6-ethylpyrimidine-4	-amine (synonym: Pylimidifen)
	Molecular formula: CAS:	C <sub>20</sub> H <sub>28</sub> ClN <sub>3</sub> O <sub>2</sub> 105779-78-0
Ò	ENCS:	Uncertain
	MW:	377.91
o cı	mp:	$70^{\circ}\text{C}^{13)}$
	bp:	Uncertain
	SW:	$2.17 \text{mg/L} (25^{\circ}\text{C})^{-13}$
/ HN— N	Specific gravity:	Uncertain
N—	logPow:	4.84 <sup>5)</sup>
[7] 1-Chloronaphthalene		
	Molecular formula:	$C_{10}H_7Cl$
ÇI	CAS:	90-13-1
	ENCS:	4-316
$\wedge$	MW:	162.62
	mp:	-2.5°C¹)
	bp:	259.3°C¹)
	SW:	0.00224% (25°C) <sup>9)</sup>
	Specific gravity:	1.19382(20/4°C) <sup>1)</sup>
V V	logPow:	3.99)
[8] 2-Chloronitrobenzene		
	Molecular formula:	C <sub>6</sub> H <sub>4</sub> ClNO <sub>2</sub>
	CAS:	88-73-3
NO <sub>2</sub>	ENCS:	3-442
	MW:	157.56
	mp:	32.5°C <sup>14)</sup>
	bp:	245.5°C <sup>14)</sup>
	SW:	441 mg/L (25°C) <sup>15)</sup>
CI	Specific gravity:	1.368 (22/4°C) <sup>16)</sup>
	logPow:	$2.24^{2)}$
[9] Salicylaldehyde		
	Molecular formula:	$C_7H_6O_2$
	CAS:	90-02-8
∠CHO	ENCS:	3-1183, 3-2660
	MW:	122.12
	mp:	-7°C <sup>9)</sup>
	bp:	197°C <sup>9)</sup>
ОН	SW:	80.8mg/L (25°C) <sup>15)</sup>
ОП	Specific gravity:	1.1674 (20/4°C) <sup>29)</sup>
	logPow:	1.81 (pH5.4) <sup>11)</sup>
[10] 2,6-Dinitrotoluene		
	Molecular formula:	$C_7H_6N_2O_4$
NO <sub>2</sub>	CAS:	606-20-2
2	ENCS:	3-446
	MW:	182.14
	mp:	64∼66°C¹ <sup>7)</sup>
ĺ Y	bp:	285°C <sup>9)</sup>
	SW:	$208 \text{ mg/L} (25^{\circ}\text{C})^{-17)}$
<u>"</u>	Specific gravity:	$1.54 \ (15/15^{\circ}C)^{-18)}$
NO <sub>2</sub>	logPow:	2.1 <sup>19)</sup>
- <b>Z</b>		

#### [11] *m*-Dinitrobenzene $C_6H_4N_2O_4$ Molecular formula: CAS: 99-65-0 ENCS: 3-445 MW: 168.11 $90.3^{\circ}C^{9)}$ mp: 291°C<sup>9)</sup> bp: SW: $2.09\%^{9)}$ NO<sub>2</sub> Specific gravity: Uncertain $1.49^{20}$ logPow: [12] Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene) Molecular formula: $C_{14}H_{14}O$ 103-50-4 CAS: 3-1082 ENCS: 198.26 MW: $3.6^{\circ}C^{9)}$ mp: 298°C<sup>9)</sup> bp: 40mg/L (35°C) <sup>21)</sup> SW: Specific gravity: 1.00142 (20/4°C) 1) 3.31<sup>11)</sup> logPow: [13] Dimethyl 4,4'-(o-phenylene)bis(3-thioallophanate) (synonym: Thiophanate-methyl) Molecular formula: $C_{12}H_{14}N_4O_4S_2\\$ CAS: 23564-05-8 ENCS: Uncertain 342.39 MW: mp: $172^{\circ}C^{9)}$ Uncertain bp: $439 mg/L (25^{\circ}C)^{-21)}$ SW: Specific gravity: Uncertain $1.5^{22}$ logPow: [14] Dimethyl terephthalate Molecular formula: $C_{10}H_{10}O_4$ 120-61-6 CAS: ENCS: 3-1328 MW: 194.19 $141^{\circ}\!\mathbb{C}^{9)}$ mp: 288°C(760 mmHg)9) bp: 19.0mg/L (25°C) <sup>21)</sup> SW: $1.065 (20/4^{\circ}\text{C})^{16)}$ $2.25^{11)}$ Specific gravity: logPow: [15] Propylene dinitrate Molecular formula: $C_3H_6N_2O_6\\$ 6423-43-4 CAS: ENCS: 2-1570 MW: 166.089 -29.49°C<sup>23)</sup> mp: O-NO2 bp: 92°C(10 mmHg)<sup>9)</sup> SW: $0.1g/100mL^{24}$ 1.2(water=1)<sup>24)</sup> 1.83<sup>25)</sup>

Specific gravity: logPow:

[16] - Niteramilia		
[16] <i>o</i> -Nitroaniline		
	Molecular formula:	$C_6H_6N_2O_2$
	CAS:	88-74-4
NO <sub>2</sub>	ENCS:	3-392
	MW:	138.13 $69 \sim 71^{\circ} \text{C}^{1)}$
	mp:	284°C <sup>1)</sup>
	bp: SW:	
NH <sub>2</sub>		1260mg/L (25°C) <sup>26)</sup>
2	Specific gravity:	0.9015 (25/4°C) <sup>1)</sup> 1.85 <sup>32)</sup>
	logPow:	1.85
[17] <i>m</i> -Nitroaniline		
	Molecular formula:	$C_6H_6N_2O_2$
	CAS:	99-09-2
	ENCS:	3-392
	MW:	138.13
	mp:	114°C¹)
	bp:	306℃¹)
O <sub>2</sub> N NH <sub>2</sub>	SW:	1g/880mL <sup>1)</sup>
$O_2N$ $\longrightarrow$ $NH_2$	Specific gravity:	0.9011 (25/4°C) <sup>1)</sup>
	logPow:	$1.37^{27}$
[18] Vanadium and its compounds (as Vanadium)		
	Molassias C 1	<b>V</b> 7
	Molecular formula: CAS:	V 7440-62-2
	ENCS:	Uncertain
	MW:	50.94
$\mathbf{V}$	mp:	1917°C¹)
▼	bp:	Uncertain
	SW:	Insoluble <sup>1)</sup>
	Specific gravity:	Uncertain
[19] Phenanthrene	logPow:	Uncertain
[17] FIICHAHUHCHC		
	Molecular formula:	$C_{14}H_{10}$
	CAS:	85-01-8
	ENCS:	4-635
	MW:	178.23
// \/ \\	mp:	$101^{\circ}C^{9}$
	bp: SW:	$340^{\circ}C^{9}$
		1.6 mg/L (15°C) <sup>17)</sup> 0.98 (4°C) <sup>9)</sup>
	Specific gravity: logPow:	$0.98 (4^{\circ}C)^{-3}$ $4.46^{2)}$
	logrow.	7.70
[20] Phenyloxirane (synonym: Styreneoxide)		
	Molecular formula:	$C_8H_8O$
O	CAS:	96-09-3
<	ENCS:	3-1033
$ \checkmark $	MW:	120.15
	mp:	-35.6°C <sup>9)</sup>
<b>/</b> /	bp:	$194.1^{\circ}C^{9}$
// \	SW:	3,000mg/L (20°C) <sup>30)</sup> 1.0523 <sup>9)</sup>
\ <u>\</u>	Specific gravity: logPow:	1.0523 <sup>37</sup> 1.61 <sup>31)</sup>
	iogi ow.	1.01

#### [21] Dimethyl phthalate Molecular formula: $C_{10} H_{10} O_4 \\$ CAS: 131-11-3 3-1301 ENCS: MW: 194.19 5.5°C¹) mp: $283.7^{\circ}C^{1)}$ bp: SW: 4g/L (20°C) <sup>18)</sup> $1.192 (20^{\circ}C)^{-16)}$ Specific gravity: 1.56<sup>32)</sup> logPow: ö [22] Benzyl alcohol C<sub>7</sub>H<sub>8</sub>O 100-51-6 Molecular formula: OH CAS: ENCS: 3-1011 MW: 108.14 $-15.2^{\circ}C^{9)}$ mp: bp: $205.3^{\circ}C^{9)}$ SW: 42.9g/L $(25^{\circ}C)^{-15)}$ 1.04 (20/4°C) 9) Specific gravity: 1.102) logPow: [23] Methylhydrazine Molecular formula: $CH_6N_2$ CAS: 60-34-4 ENCS: 2-2385 MW: 46.07 $-52.4^{\circ}C^{1)}$ mp: $NH-NH_2$ bp: $87.5^{\circ}C^{1)}$ SW: $1,000 \text{ g/L} (25^{\circ}\text{C})^{-33)}$ Specific gravity: $0.874~(25^{\circ}C)^{-1)}$ $-1.05^{34}$ logPow: [24] 2-Methyl-1,1'-biphenyl-3-ylmethyl (Z)-3-(2-chloro-3,3,3-trifluoro-1-propenyl)-2,2-dimethylcyclopropanecarboxylate (synonym: Bifenthrin) Molecular formula: $C_{23}H_{22}ClF_3O_2\\$ CAS: 82657-04-3 ENCS: 4-1701 422.87 MW: 69°C1) mp: H Uncertain bp: $0.1 mg/L^{13)} \\$ SW: Specific gravity: 1.21 (25°C) 1) logPow: >6<sup>2)</sup>



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# 3. Surveyed site and procedure

In the Initial Environmental Survey, the sampling and analysis of specimens was entrusted to prefectural governments and government-designated cities across Japan, and some specimens were analyzed by private analytical laboratories.

## (1) Organisations responsible for sampling

Local		Survey	yed media	
communities	Organisations responsible for sampling	Surface water	Sediment	Air
Hokkaido	Hokkaido Institute of Environmental Sciences	0	0	0
Sapporo City	Sapporo City Institute of Public Health	0		0
Iwate Pref.	Research Institute for Environmental Sciences and Public Health of Iwate Prefecture	0		
Miyagi Pref.	Miyagi Prefectural Institute of Public Health and Environment	0		
Sendai City	Sendai City Institute of Public Health	0		0
Yamagata Pref.	Environmental Science Research Center of Yamagata Prefecture	0		
Ibaraki Pref.	Ibaraki Kasumigaura Environmental Science Center	0		0
Tochigi Pref.	Tochigi Prefectural Institute of Public Health and Environmental Science	0		
Saitama Pref.	Center for Environmental Science in Saitama	0		0
Chiba Pref.	Chiba Prefectural Environmental Research Center	0		0
Tokyo Met.	Tokyo Metropolitan Research Institute for Environmental Protection			0
Kanagawa	Kanagawa Environmental Research Center			
Pref.	-			0
Yokohama	Yokohama Environmental Science Research Institute	0		
City				
Kawasaki	Kawasaki Municipal Research Institute for Environmental Protection	0	0	0
City				
Niigata Pref.	Niigata Prefectural Institute of Public Health and Environmental Sciences	0	0	
Ishikawa Pref.	Ishikawa Prefectural Institute of Public Health and Environmental Science	0	0	
Nagano Pref.	Nagano Environmental Conservation Research Institute	0		0
Gifu Pref.	Gifu Prefectural Research Institute for Health and Environmental Sciences			0
Aichi Pref.	Aichi Environmental Research Center	0	0	
Nagoya City	Nagoya City Environmental Science Research Institute	0		0
Mie Pref.	Mie Prefectural Science and Technology Promotion Center	0	0	0
Shiga Pref.	Lake Biwa Environmental Research Institute	0	0	
Kyoto Pref.	Kyoto Prefectural Institute of Public Health and Environment	0		0
Kyoto City	Kyoto Prefectural Institute of Public Health and Environment	0		0
Osaka Pref.	Research Institute of Environment, Agriculture and Fisheries, Osaka Prefectural Government	0		0
Osaka City	Osaka City Institute of Public Health and Environmental Sciences	0		
Hyogo Pref.	Hyogo Prefectural Institute of Public Health and Environmental Sciences	0	0	
Wakayama Pref.	Wakayama Prefectural Research Center of Environment and Public Health			0
Okayama Pref.	Okayama Prefectural Institute for Environmental Science and Public Health	0	0	0
Hiroshima City	Hiroshima Prefectural Technology Research Institute Health and Environment Center	0	0	
Yamaguchi Pref.	Yamaguchi Prefectural Public Health and Environment	0	0	0
Tokushima Pref.	Tokushima Prefectural Institute of Public Health and Environmental Sciences	0		0
Kagawa Pref.	Kagawa Prefectural Research Institute for Environmental Sciences and Public Health	0	0	0
Fukuoka Pref.	Fukuoka Institute of Health and Environmental Science	0		0

Local	Organizations resmansible for semuling	Survey	ved media	
communities	Organisations responsible for sampling	Surface water	Sediment	Air
Kitakyushu	Kitakyushu City Institute of Environmental Sciences		0	
City		O		
Fukuoka	Fukuoka City Institute for Hygiene and the Environment			
City		O		
Saga Pref.	Saga Prefectural Environmental Research Center	0	0	0
Kumamoto	Kumamoto Prefectural Institute of Public Health and Environmental			)
Pref.	Science			

(Note) Organisations responsible for sampling are described by their official names in FY 2007.

## (2) Surveyed sites and target chemicals

Surveyed sites and target chemicals for surface water are shown in Table 1-1-1 and Figure 1-1-1. Surveyed sites and target chemicals for sediment are shown in Table 1-1-2 and Figure 1-1-1. Surveyed sites and target chemicals for air are shown in Table 1-1-3 and Figure 1-1-2. The breakdown is summarized as follows.

Surveyed media	Numbers of local communities	Numbers of target chemicals	Numbers of surveyed sites	Numbers of samples at a surveyed site
Surface water	33	17	53	3
Sediment	14	3	18	3
Air	24	10	29	3

Table 1-1-1 List of surveyed sites (surface water) and target chemicals in the Initial Environmental Survey in FY 2007

2007	1								,									
Local communities	Surveyed sites	F 4 7	re1	TO 3	TO1	F1.07	F1 1 7		arget				F107	[20]	[21]	[22]	F2 41	[27]
Hokkaido	Ishikarikakokyo Bridge, Mouth of Riv.Ishikari(Ishikari City)	[4]	[5]	[8]	[9]	[10]	[11]	$\bigcirc$	[13]	[16]	[17]			[20]	[21]	[23]	[24]	[26]
Sapporo City	Nakanuma of Riv. Toyohira(Sapporo City)																	
Support City	Azuma-bashi Bridge, Riv.Toyohira(Sapporo City)							0									0	
Iwate Pref.	Riv.Toyosawa(Hanamaki City)											0					)	
Miyagi Pref.	Kannon-Bridge of Riv.Naruse (Misato Town)								0									
Sendai City	Hirose-ohashi Bridge, Riv.Hirose(Sendai City)					0	0			0	0							
Yamagata Pref.	Mouth of Riv.Mogami(Sakata City)															0		
Ibaraki Pref.	Katta-bashi Bridge, Riv.Naka(Hitachinaka City)					0	0		0	0	0						0	
	Tonekamome-ohasi Bridge, Mouth of Riv.Tone(Kamisu City)					0	0		0	0	0						0	
Tochigi Pref.	Riv.Tagawa (Utsunomiya City)	0															0	
Saitama Pref.	Nawate-Bridge of Riv.Ayase(Saitama City)				0													
Chiba Pref.	Asahi Higashiashiarai Beach													0				
	Asai-bashi Bridge, Riv.Yourou(Ichihara City)								0							0		
Yokohama City	Kamenoko Bridge over Riv.Tsurumi(Yokohama City)								0			0	0				0	
Kawasaki City	Mouth of Riv. Tama(Kawasaki City)			0		0	0	0		0	0				0			
	Keihin Canal, Port of Kawasaki			0		0	0	0		0	0				0			
Niigata Pref.	Lower Riv. Shinano(Niigata City)		0	0		0	0	0		0	0				0			
Ishikawa Pref.	Mouth of Riv. Sai(Kanazawa City)	0	0	0		0	0	0		0	0							
	Offshore of Kasama Town(Hakusan City)																	0
Nagano Pref.	Lake Suwa (center)	0	0														0	
Aichi Pref.	Nagoya Port												0					
Nagoya City	Minatoshinbashi Bridge, Riv. Hori (Nagoya City)															0		
Mie Pref.	Yokkaichi Port	0	0	0		0	0			0	0		0		0			
	Tsumatsusaka Port																	0
Shiga Pref.	Nomura-Bridge of Riv. Hino(Omihachiman City) Hattori-Bridge of Riv. Yasu(Moriyama City)				0									0				
	Lake Biwa(center, offshore of Minamihira)		0										0					
Kyoto Pref.	Miyazu Port							0				0						
Kyoto City	Miyamae Bridge, Riv. Katsura(Kyoto City)															0		
Osaka Pref.	Dainineyagawa, Shinkingoro Bridge(Higashiosaka City)																	0
	Mouth of Riv. Yamato (Sakai City)		0													0		
Osaka City	Osaka Port				0													
Hyogo Pref.	Amagasaki-Nishinomiya-Ashiya Port Awata-Bridge of Riv. Kako(Ono City)													0				0
	Offshore of Himeji		0					0										
Okayama Pref.	Offshore of Mizushima							Ī	0				0		0		0	
Hiroshima Pref.	Mitsu Bay				0								Ė					
	Kure Port														0			
	Hiroshima Bay														0			
	West of Hiroshima Bay													0				

Local	Surveyed sites	Target chemicals								•								
communities	Surveyed sites	[4]	[5]	[8]	[9]	[10]	[11]	[12]	[13]	[16]	[17]	[18]	[19]	[20]	[21]	[23]	[24]	[26]
Yamaguchi Pref.	East of Tokuyama Bay																	$\circ$
	Tokuyama Bay												0				$\circ$	
	Offshore of Ube				0													
	Offshore of Hagi												0				$\circ$	
Tokushima Pref.	Mouth of Riv. Yoshino(Tokushima City)								0									
Kagawa Pref.	Takamatsu Port	0											0				0	
Fukuoka Pref.	Offshore of Omuta	0		0					0									
	Kabura-bashi Bridge, River Raizan(Maebaru City)	0							0									
Kitakyushu City	Kanmon Strait	0		0		0	0			0	0		0					
	Dokai Bay	0		0		0	0			$\bigcirc$	0		0					
Fukuoka City	Hakata Bay											0						
Saga Pref.	Mameda-Bridge of Riv. Iryuu(Yoshinogari Town)													0				
	Imari Bay	0						0										

<sup>[4] 2,4-</sup>Xylenol, [5] Quinoline, [8] 2-Chloronitrobenzene, [9] Salicylaldehyde, [10] 2,6-Dinitrotoluene, [11] *m*-Dinitrobenzene, [12] Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene), [13] Dimethyl 4,4'-(o-phenylene)bis(3-thioallophanate) (synonym: Thiophanate-methyl), [16] *o*-Nitroaniline, [17] *m*-Nitroaniline, [18] Vanadium and its compounds (as Vanadium), [19] Phenanthrene, [20] Phenyloxirane (synonym: Styreneoxide), [21] Dimethyl phthalate, [23] Methylhydrazine, [24] 2-Methyl-1,1'-biphenyl-3-ylmethyl (Z)-3-(2-chloro-3,3,3-trifluoro-1-propenyl)-2,2-dimethylcyclopropanecarboxylate (synonym: Bifenthrin), [26] Mercaptoaceticacid

Table 1-1-2 List of surveyed sites (sediment) and target chemicals in the Initial Environmental Survey in FY 2007

I and annumities	Comment sites	Targe	et chen	nicals
Local communities	Surveyed sites	[12]	[19]	[21]
Hokkaido	Ishikarikakokyo Bridge, Mouth of Riv. Ishikari(Ishikari City)	0	0	
Kawasaki City	Mouth of Riv. Tama(Kawasaki City)	0		0
	Keihin Canal, Port of Kawasaki	0		0
Niigata Pref.	Lower Riv. Shinano(Niigata City)	0		0
Ishikawa Pref.	Mouth of Riv. Sai(Kanazawa City)	0		
Aichi Pref.	Nagoya Port		0	
Mie Pref.	Yokkaichi Port		0	0
Shiga Pref.	Lake Biwa(center, offshore of Minamihira)		0	
Hyogo Pref.	Offshore of Himeji	0		
Okayama Pref.	Offshore of Mizushima		0	
Hiroshima Pref.	Kure Port			0
	Hiroshima Bay			0
Yamaguchi Pref.	Tokuyama Bay		0	
	Offshore of Hagi		0	
Kagawa Pref.	Takamatsu Port		0	
Kitakyushu City	Kanmon Strait		0	
	Dokai Bay		0	
Saga Pref.	Imari Bay	0		

<sup>[12]</sup> Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene), [19] Phenanthrene, [21] Dimethyl phthalate

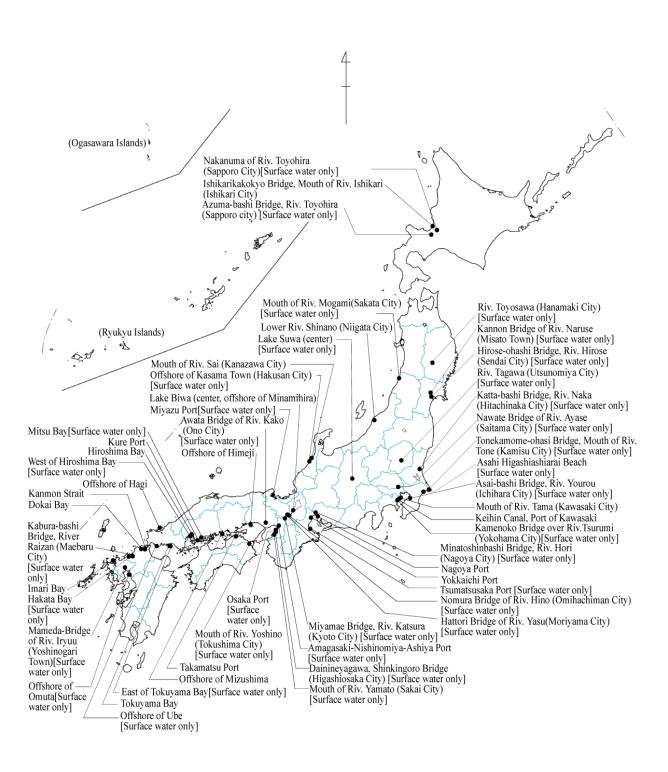


Figure 1-1-1 Surveyed sites (surface water and sediment) in the Initial Environmental Survey in FY 2007

Table 1-1-3 List of surveyed sites (air) and target chemicals in the Initial Environmental Survey in FY 2007

1 4010 1-1	3 List of surveyed sites (air) and target chemicals in the Initial Environn	Litte	ii Su	nvey			chem	icals			
Local communities	Surveyed sites	[1]	[2]	[3]	[6]	[7]	[14]	[15]	[22]	[25]	[27]
Hokkaido	Hokkaido Institute of Environmental Sciences(Sapporo City)	0				0	0			0	
Sapporo City	Sapporo City Institute of Public Health(Sapporo City)								0		0
Sendai City	Tsutsujigaoka Park(Sendai City)						0		0		
Ibaraki Pref.	Ibaraki Kasumigaura Environmental Science Center(Tsuchiura City)			0	0					0	
Saitama Pref.	Center for Environmental Science in Saitama(Kisai Town)						0	0			0
Chiba Pref.	Chiba Prefectural Environmental Research Center(Ichihara City)		0	0	0		0			0	0
	Ichihara-Matsuzaki Air Quality Monitoring Station(Ichihara City)					0					
Tokyo Met.	Tokyo Metropolitan Research Institute for Environmental Protection(Koto Ward)					0					
	Chichijima Island					0					
Kanagawa Pref.	Sagamihara City Fuchinobe Elementary School(Sagamihara City)		0	0						0	
	Kanagawa Environmental Research Center(Hiratsuka City)	0									0
Kawasaki City	Daishi Air Quality Monitoring Station(Kawasaki City)		0	0						0	
Nagano Pref.	Nagano Environmental Conservation Research Institute(Nagano City)					0	0				
Gifu Pref.	Gifu Prefectural Research Institute for Health and Environmental Sciences(Kakamigahara City)							0			
Nagoya City	Chikusa Ward Heiwa Park(Nagoya City)	0		0	0		0	0			
Mie Pref.	Mie Prefectural Science and TechnologyPromotion Center(Yokkaichi City)					0	0	0	0		
Kyoto Pref.	Kyoto Prefectural Joyo High School(Joyo City)					0					
Kyoto City	Kyoto City Hall(Kyoto City)	0							0		
Osaka Pref.	Research Institute of Environment, Agriculture and Fisheries, Osaka Prefectural Government(Osaka City)					0	0				
Wakayama Pref.	Wakayama Prefectural Research Center of Environment and Public Health(Wakayama City)	0									
Okayama Pref.	Shionasu Air Quality Monitoring Station(Kurashiki City)		0	0						0	
Yamaguchi Pref.	Shunan City Government Building(Shunan City)		0	0						0	
	Yamaguchi Prefectural Public Health and Environment(Yamaguchi City)						0	0	0		0
Tokushima Pref.	Wakimachi ambient air pollution monitoring station(Mima City)		0	0						0	
Kagawa Pref.	Takamatsu Joint Prefectural Government Building(Takamatsu City)							0	0		
Fukuoka Pref.	Omuta City Hall(Omuta City)							0			
	Munakata Prefectural Government Building(Munakata City)							0			
Saga Pref.	Saga Prefectural Environmental Research Center(Saga City)			0	0						
Kumamoto Pref.	Kumamoto Prefectural Institute of Public Health and Environmental Science(Udo City)				0		0				

<sup>[1]</sup> Adipic acid, [2] Ethyleneimine, [3] 4'-Ethoxyacetanilide (synonym:Phenacetin), [6] 5-Chloro-*N*-{2-[4-(2-ethoxyethyl)-2,3-dimethylphenoxy]ethyl}-6-ethylpyrimidine-4-amine (synonym:Pylimidifen), [7] 1-Chloronaphthalene, [14] Dimethyl terephthalate, [15] Propylene dinitrate, [22] Benzyl alcohol,

<sup>[25] 2-(1-</sup>Methylpropyl)-4,6-dinitrophenol, [27] Triphenyl phosphate

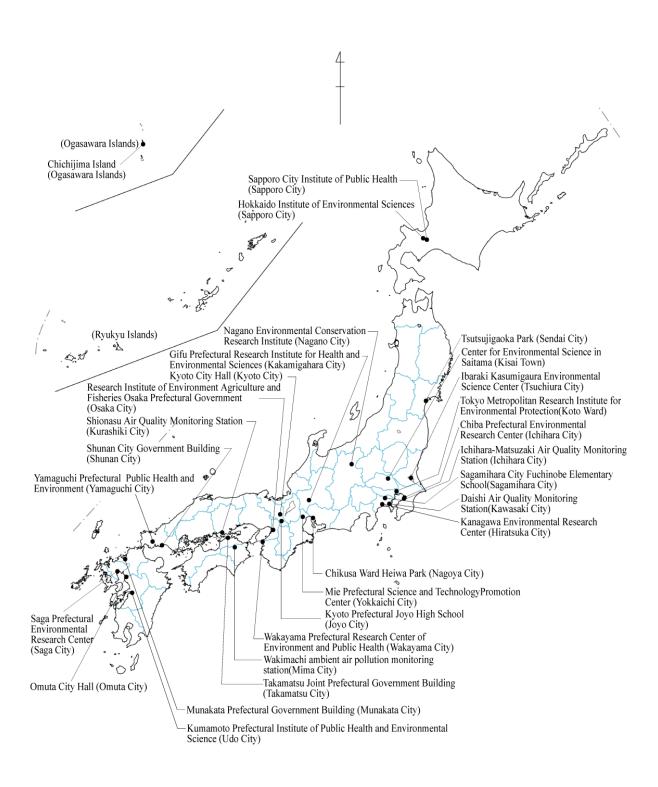


Figure 1-1-2 Surveyed sites (air) in the Initial Environmental Survey in FY 2007

### (3) Detection limit

The detection limits of analysed values reported by the analytical laboratory are not necessarily the same because of differences in the properties of specimens and in the available measurement equipment. To enable summarisation, therefore, a unified detection limit is predetermined and the analytical values reported by the analytical laboratory are summarised by the following procedure.

Treatment of measured value as an undetected value in high-sensitivity analysis

In the case of high-sensitivity analysis, in which the detection limit of the analytical laboratory is lower than the unified detection limit, any measured value lower than the unified detection limit is treated as an undetected value in the nationwide summary (see schematic ①).

Elimination of undetected values in low-sensitivity analysis from summary subject. When the detection limit of the analytical laboratory is higher than the unified detection limit, any target chemical not detected is eliminated from the subject of the summary (see schematic ②).

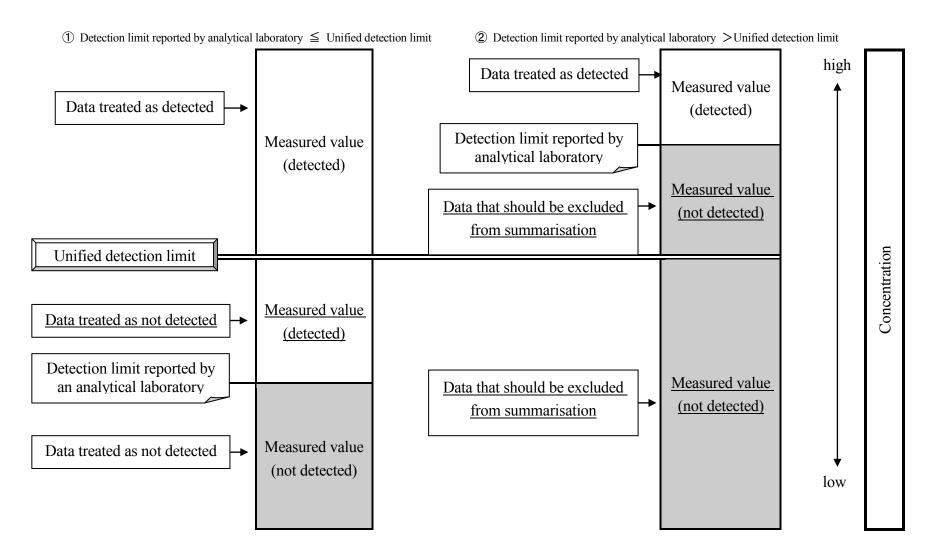
When the instrument detection limit (IDL) and the method detection limit (MDL) are given in the analytical method, which is described in reports on the investigation of the development of analytical methods for chemicals and adopted in the Initial Environmental Survey (hereafter, the Initial Environmental Survey Analytical Method), if the IDL measured by the analytical laboratory is lower than the given IDL, the MDL of the Initial Environmental Survey Analytical Method is used as the detection limit by the analytical laboratory.

When IDL and MDL are not given in the Initial Environmental Survey Analytical Method, the detection limit is predetermined by the following procedure.

When the analytical laboratory calculates the appropriate IDL and MDL following the calculation method stated in the analytical method development instruction manuals, this calculated MDL is used as the detection limit by the analytical laboratory.

When the appropriate IDL and MDL are not calculated by the analytical laboratory, one of the following procedures was employed to establish the detection limit by the analytical laboratory.

- deduction from the IDL and MDL calculated for the corresponding chemical by Initial Environmental Survey Analytical Method or other analytical laboratories
- deduction from the lowest calibration curve concentration and the results of recovery tests
- deduction from the results of addition and collection tests, the results of operation blank tests, and the signal/noise ratio (S/N ratio) obtained from the chromatogram of environmental specimens



Schematic of procedure for data summarisation

## 4. Summary of survey results

The detection ranges and the detection limits are shown in Table 1-2. The survey results are summarized as follows.

In surface water, 8 out of the 17 target chemicals were detected.

- •[4] 2,4-Xylenol: 5 of the 9 valid sites
- •[5] Quinoline: 4 of the 7 valid sites
- •[12] Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene): 1 of the 8 valid sites
- •[13] Dimethyl 4,4'-(o-phenylene)bis(3-thioallophanate) (synonym: Thiophanate-methyl): 1 of the 9 valid sites
- •[18] Vanadium and its compounds (as Vanadium): 5 of the 5 valid sites
- •[19] Phenanthrene: 6 of the 9 valid sites
- •[21] Dimethyl phthalate: 7 of the 7 valid sites
- •[26] Mercaptoaceticacid: 3 of the 5 valid sites

In sediment, 3 out of the 3 target chemicals were detected.

- •[12] Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene): 3 of the 6 valid sites
- •[19] Phenanthrene: 10 of the 10 valid sites
- •[21] Dimethyl phthalate: 6 of the 6 valid sites

In air, 5 out of the 10 target chemicals were detected.

- •[7] 1-Chloronaphthalene: 5 of the 8 valid sites
- •[14] Dimethyl terephthalate: 8 of the 9 valid sites
- •[15] Propylene dinitrate: 1 of the 8 valid sites
- •[22] Benzyl alcohol: 5 of the 6 valid sites
- •[27] Triphenyl phosphate: 5 of the 5 valid sites

Table 1-2 Summary of the detection ranges and the detection limits in the Initial Environmental Survey in FY 2007

Table 1-2	2 Summary of the detection ranges and	Surface wa		Sediment [		Air [ng	
	Target chemicals	Detection Detection	ter [ng/L]	Detection			ym j
No.	Detection range and frequency	range and frequency	Detection limit	range and frequency	Detection limit	Detection range and frequency	Detection limit
[1]	Adipic acid					nd 0/5	90
[2]	Ethyleneimine					nd 0/6	2.7
[3]	4'-Ethoxyacetanilide (synonym: Phenacetin)					nd 0/9	3.1
[4]	2,4-Xylenol	nd~4.3 5/9	1.4				
[5]	Quinoline	nd~13 4/7	1.1				
[6]	5-Chloro- <i>N</i> -{2-[4-(2-ethoxyethyl)-2,3-di methylphenoxy]ethyl}-6-ethylpyrimidin e-4-amine (synonym: Pylimidifen)					nd 0/5	1.8
[7]	1-Chloronaphthalene					nd∼0.73 5/8	0.15
[8]	2-Chloronitrobenzene	nd 0/8	2.3				
[9]	Salicylaldehyde	nd 0/5	13				
[10]	2,6-Dinitrotoluene	nd 0/7	1.4				
[11]	<i>m</i> -Dinitrobenzene	nd 0/8	1.9				
[12]	Dibenzyl ether (synonym: [(Benzyloxy)methyl]benzene)	nd~8.3 1/8	1.9	nd~21 3/6	0.18		
[13]	Dimethyl 4,4'-(o-phenylene)bis(3-thioallophanate) (synonym: Thiophanate-methyl)	nd~0.90 1/9	0.79				
[14]	Dimethyl terephthalate					nd~1.0 8/9	0.012
[15]	Propylene dinitrate					nd~3.9 1/8	2.0
[16]	o-Nitroaniline	nd 0/8	1.1				
[17]	<i>m</i> -Nitroaniline	nd 0/7	2.2				
[18]	Vanadium and its compounds	630~4,600 5/5	2.4				
[19]	Phenanthrene (as Vanadium)	nd∼55 6/9	1.4	3.9~690 10/10	0.023		
[20]	Phenyloxirane (synonym: Styreneoxide)	nd 0/5	12				
[21]	Dimethyl phthalate	nd~9.7 7/7	1.7	0.54~6.3 6/6	0.35		
[22]	Benzyl alcohol					nd~7,300 5/6	450
[23]	Methylhydrazine	nd 0/5	27				
[24]	2-Methyl-1,1'-biphenyl-3-ylmethyl( <i>Z</i> )-3 -(2-chloro-3,3,3-trifluoro-1-propenyl)-2, 2-dimethylcyclopropanecarboxylate (synonym: Bifenthrin)	nd 0/11	7.8				
[25]	2-(1-Methylpropyl)-4,6-dinitrophenol					nd 0/8	3.2
[26]	Mercaptoaceticacid	nd∼24 3/5	1.1				
[27]	Triphenyl phosphate					0.054~0.33 5/5	0.041

(Note 1) Detection frequency is based on the number of sites or areas, thus means (the number of detected sites/the number of surveyed sites). A site where data was not available was excluded from the number of surveyed sites. A site where the data became invalid under a unified detection limit was also excluded. 3 samples were measured for a site or area, and the detection in more than one out of samples from a site or area can be defined as one detected site or area. (Note 2) Detection range is based on the number of samples and therefore can be shown as "nd~" even if a target chemical is detected in all sites (or areas). (Note 3) means the medium was not surveyed.