# Chapter 2 Results of the Detailed Environmental Survey in FY 2006

# 1. Purpose of the survey

The Detailed Environmental Survey is aimed at understanding the environmental persistence of the Specified Chemical Substances and Monitored Chemical Substances under the Law Concerning the Examination and Regulation of Manufacture, etc. of Chemical Substances (Law No. 117 of 1973) (hereafter, the Chemical Substances Control Law) and chemicals requiring the Initial Environmental Risk Assessment.

# 2. Target chemicals

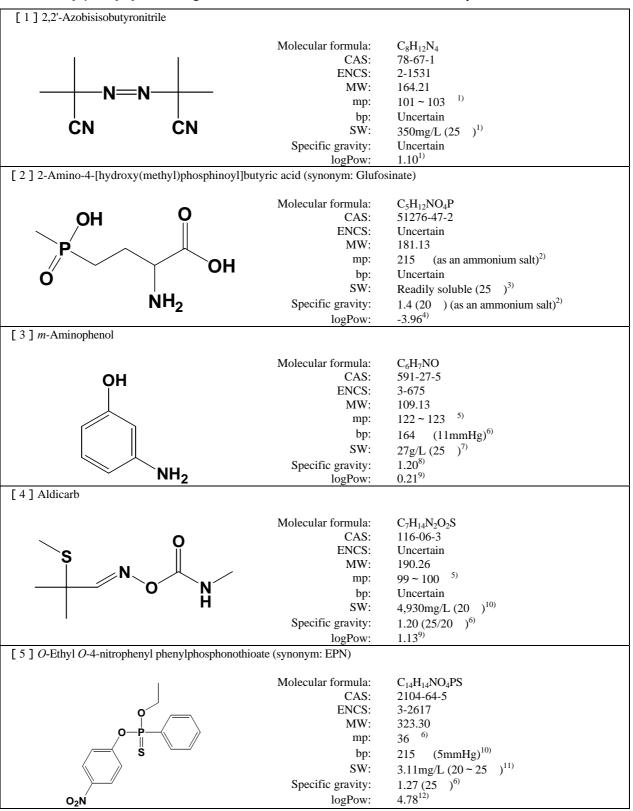
In the FY 2006 Detailed Environmental Survey, 38 chemicals (groups) that were selected and designated as target chemicals. The combinations of target chemicals and surveyed media are given below.

	Target chemicals	Designated	Class in		Survey	ed media		
		The Chemical	The PRTR	Surface				
No	Name	Substances	Law	water	Sediment	Wildlife	Air	food
[1]	2,2'-Azobisisobutyronitrile	Control Law II Monitored	I					
	2-Amino-4-[hydroxy(methyl)phosphinoyl]but							
[2]	yric acid (synonym: Glufosinate)		I					
[3]	m-Aminophenol	II Monitored	I					
		III Monitored	1					
[4]	Aldicarb							
[5]	O-Ethyl O-4-nitrophenyl	II Monitored	I					
	phenylphosphonothioate (synonym: EPN)	III Monitored						
	<i>N,N'</i> -Ethylenebis(dithiocarbamamic acid) and							
	its salt (as a common component of originally							
	targeted chemicals [6-1] ~ [6-3])			ļ			ļ 	
	[6-1] Manganese <i>N,N'</i> -ethylenebis(dithiocarbamate) (synonym:		I					
	Maneb)		1					
[6]	[6-2] Complex compounds of manganese			1		<b></b>		
	N,N'-ethylenebis(dithiocarbamate) and zinc		т					
	<i>N,N'</i> -ethylenebis(dithiocarbamate) (synonym:		I					
	Mancozeb)			ļ				
	[6-3] <i>N,N'</i> -Ethylenebis(thiocarbamoylthiozinc)		<b>.</b>					
	bis( <i>N</i> , <i>N</i> -dimethyldithiocarbamate) (synonym: Polycarbamate)		I					
[7]	2,6-Xylenol	III Monitored	I					
[8]	Chlorobenzene	III Monitored	I					
[9]	Isobutyl acetate							
[10]	Diisopropylnaphthalene	I Monitored						
	S-4-Chlorobenzyl N,N-diethylthiocarbamate		T					
[11]	(synonym: Thiobencarb)		I					
[12]	3-(3,4-Dichlorophenyl)-1,1-dimethylurea	II Monitored	I					
[12]	(synonym: Diuron or DCMU)	11 Womtored	•					
[13]	2,6-Dichlorobenzonitrile (synonym: Dichlobenil)	II Monitored	I					
	2,4-Di- <i>tert</i> -butyl-6-(5-chloro-2 <i>H</i> -1,2,3-							
[14]	benzotriazol-2-yl)phenol	I Monitored						
	<i>N,N</i> -Dimethyldithiocarbamamic acid and its							
	salt (as a common component of originally							
	targeted chemicals [15-1] ~ [15-2])							
[15]	[15-1] Zinc bis( <i>N</i> , <i>N</i> -dimethyldithiocarbamate)	II Monitored	J					
	(synonym: Ziram)		*	ļ		}	 	
	[15-2] <i>N,N'</i> -Ethylenebis(thiocarbamoylthiozinc) bis( <i>N,N</i> -dimethyldithiocarbamate) (synonym:		I	1				
	Polycarbamate)		1	1				
[16]	N,N-Dimethyldodecylamine N-oxide		I	1				
[17]	<i>N,N</i> -Dimethylformamide	II Monitored	I					
[18]	Hydrogenated terphenyl	I Monitored	-					
[10]	, 50	1 111011110101	l .	1	l	l	l	

	Target chemicals	Designated (	Class in		Survey	ed media		
No	Name	The Chemical Substances Control Law	The PRTR Law	Surface water	Sediment	Wildlife	Air	food
[19]	O,O-Diethyl O-2-isopropyl-6-methyl-4-pyrimidinyl phosphorothioate (synonym: Diazinon)	II Monitored III Monitored	I					
[20]	O,O-Dimethyl O-3-methyl-4-nitrophenyl phosphorothioate (synonym: Fenitrothion or MEP)	II Monitored	I					
[21]	Tetrachloroisophthalonitrile (synonym: Chlorothalonil or TPN)	II Monitored	I					
[22]	Tetrahydrofuran							
[23]	Trichloroacetaldehyde	II Monitored	I					
[24]	Trichloronitromethane (synonym: Chloropicrin)	II Monitored	I					
[25]	Nitrofen (synonym: NIP)							
[26]	1,1-Bis( <i>tert</i> -butyldioxy)-3,3,5- trimethylcyclohexane	I Monitored						
[27]	Hydrazine	II Monitored III Monitored	I					
[28]	1-Butanol							
[29]	Furfural							
[30]	2-(2-Benzothiazolyloxy)- <i>N</i> -methylacetanilide (synonym: Mefenacet)		I					
[31]	2-(2 <i>H</i> -1,2,3-Benzotriazol-2-yl)-4,6-di- <i>tert</i> -but ylphenol	I Specified						
[32]	Methyl methacrylate		I					
[33]	2-(1-Methylethoxy)ethanol							
[34]	2,3-Dihydro-2,2-dimethyl-7-benzo[b]furanyl <i>N</i> -methylcarbamate (synonym: Carbofuran)	II Monitored III Monitored	I					
[35]	2-sec-Butylphenyl <i>N</i> -methylcarbamate (synonym: Fenobucarb or BPMC)	II Monitored III Monitored	I					
[36]	α-Methylstyrene	III Monitored	I					
[37]	Dimethyl 2,2-dichlorovinyl phosphate (synonym: Dichlorvos or DDVP)	II Monitored III Monitored	I					
[38]	Tributyl phosphate	II Monitored	I (Note 2)					

(Note 1) "The PRTR Law" hereafter means "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)." (Note 2) Only as tri-*n*-butyl phosphate is designated.

Chemical and physical properties of target chemicals of the Detailed Environmental Survey are as follows.



(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.

```
[ 6 ] N,N'-Ethylenebis(dithiocarbamamic acid) and its salt
                                                        Molecular formula:
                                                                                 C_4H_8N_2S_4
                                                                      CAS:
                                                                                 115-54-6 etc.
                                                                                 2-1808
                                                                     ENCS:
                                                                                 212.38
                                                                      MW:
                                SH
                     NH
                                                                        mp:
                                                                                 Uncertain
                                                                                 Uncertain
                                                                        bp:
                     NH
                                SH
                                                                       SW:
                                                                                 114g/L (25)^{3}
                                                           Specific gravity:
                                                                                 Uncertain
                                                                   logPow:
                                                                                 0.62^{4)}
[ 6-1 ] Manganese N,N'-ethylenebis(dithiocarbamate) (synonym: Maneb)
                                                        Molecular formula:
                                                                                 C<sub>4</sub>H<sub>6</sub>MnN<sub>2</sub>S<sub>4</sub>
                                                                                 12427-38-2
                                                                      CAS:
                                                                     ENCS:
                                                                                 2-1841
                                                                      MW:
                                                                                 289.32
                                                                                               (decomposition)<sup>10)</sup>
                     ·NH
                                                                        mp:
                                                                                 192 ~ 204
                                                                        bp:
                                                                                 Uncertain
                     NΗ
                                                                       SW:
                                                                                 6mg/L (25
                                                           Specific gravity:
                                                                                 1.92 (25/4
                                                                                 0.62^{4)}
                                                                   logPow:
[ 6-2 ] Complex compounds of manganese N,N'-ethylenebis(dithiocarbamate) and zinc N,N'-ethylenebis(dithiocarbamate)
        (synonym: Mancozeb)
                                                        Molecular formula:
                                                                                 C_4H_6MnN_2S_4/C_4H_6N_2S_4Zn
                                                                      CAS:
                                                                                 8018-01-7
                                                                     ENCS:
                                                                                 2-2127
                                                                      MW:
                                                                                 289.32/299.79
                                                                                               (decomposition)<sup>10)</sup>
                                                                        mp:
                                                                                 192 ~ 204
     -NH
                                                                        bp:
                                                                                 Uncertain
                                                                       SW:
                                                                                 6.2mg/L (25
                                                           Specific gravity:
                                                                                 Uncertain
                                                                                 1.33^{10)}
                                     S
                                                                  logPow:
[ 6-3 ] N,N'-Ethylenebis(thiocarbamoylthiozinc) bis(N,N-dimethyldithiocarbamate) (synonym: Polycarbamate)
                                                        Molecular formula:
                                                                                 C_{10}H_{18}N_4S_8Zn_2\\
                                                                      CAS:
                                                                                 64440-88-6
                                                                     ENCS:
                                                                                 2-1848
                                                                      MW:
                                                                                 581.61
                                                                        mp:
                                                                                 Uncertain
                                                                                 Uncertain
                                                                        bp:
                                                                       SW:
                                                                                 Uncertain
                                                           Specific gravity:
                                                                                 Uncertain
                                                                   logPow:
                                                                                 Uncertain
[7] 2,6-Xylenol
                                                        Molecular formula:
                                                                                 C_8H_{10}O
                                                                                 576-26-1
                                                                      CAS:
                         OH
                                                                     ENCS:
                                                                                 3-521, 4-57
                                                                      MW:
                                                                                 122.16
                                                                                 49 5)
                                                                        mp:
                                                                                 203 5)
                                                                        bp:
                                                                       SW:
                                                                                 6,050 \text{mg/L} (25)^{7}
                                                                                 1.13^{5)}
                                                           Specific gravity:
                                                                                 2.36^{9)}
                                                                   logPow:
```

[8] Chlorobenzene	
	G M GI
Molecular formula:	C <sub>6</sub> H <sub>5</sub> Cl
CAS: ENCS:	108-90-7 3-31
MW:	112.56
mp:	-45.2 <sup>6)</sup>
bp:	131.7 6)
SW:	498mg/L (25 ) <sup>14)</sup>
Specific gravity:	1.11 (20 ) <sup>6)</sup>
logPow:	2.89 <sup>9)</sup>
[ 9 ] Isobutyl acetate	2.07
[ 7 ] Isobuty Luceture	
Molecular formula:	$C_6H_{12}O_2$
CAS:	110-19-0
ENCS:	2-731
MW:	116.16
mp:	-98.8 <sup>6)</sup>
bp:	116.5 6)
SW:	$6,300 \text{mg/L} (25)^{7}$
Specific gravity:	$0.871 (20/4)^{6}$
logPow:	1.789)
[ 10 ] Diisopropylnaphthalene	
Molecular formula:	$C_{16}H_{20}$
CAS:	38640-62-9
ENCS:	4-961
MW:	212.33
mp: bp:	Uncertain 290 ~ 299 1)
SW:	$0.11 \text{mg/L} (25)^{15}$
Specific gravity:	0.96 (25 ) <sup>6</sup>
logPow:	$6.08^{4}$
[ 11 ] S-4-Chlorobenzyl N,N-diethylthiocarbamate (synonym: Thiobencarb)	0.00
[ 11 ] 5-4-Chiotoochizyi 14,14-dichiyidhocaroaniate (synonyin, 11hoochicato)	
Molecular formula:	$C_{12}H_{16}CINOS$
CAS:	28249-77-6
ENCS:	Uncertain
MW:	257.78
mp:	3.3 10)
op:	$126 \sim 129 \qquad (0.0008 \text{mmHg})^6$
SW:	28mg/L (25 ) <sup>13)</sup>
Specific gravity:	$1.15 \sim 1.18 (20)^{10}$
CI logPow:	3.40 <sup>9)</sup>
[ 12 ] 3-(3,4-Dichlorophenyl)-1,1-dimethylurea (synonym: Diuron or DCMU	)
Molecular formula:	$C_9H_{10}Cl_2N_2O$
CAS:	330-54-1
HN N ENCS:	3-2194
MW:	233.09 158 ~ 159 5)
mp:	130 137
bp:	$180 \sim 190$ (decomposition) <sup>2)</sup> $36.4 \text{mg/L} (25)^{10}$
CI SW:	2 \ ,
Specific gravity: CI Specific gravity:	Uncertain 2.68 <sup>9)</sup>
logPow:	2.00

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[ 13 ] 2,6-Dichlorobenzonitrile (synonym: Dichlobenil)
                                                      Molecular formula:
                                                                              C_7H_3Cl_2N
                                                                    CAS:
                                                                               1194-65-6
                         CN
                                                                  ENCS:
                                                                               3-4103
                                                                    MW:
                                                                               172.01
                                                                               144 ~ 145
                                                                     mp:
            CI
                                    CI
                                                                              270^{-10)}
                                                                     bp:
                                                                               14.6mg/L (20 )<sup>10)</sup>
                                                                     SW:
                                                                              Uncertain
                                                         Specific gravity:
                                                                              2.74^{9)}
                                                                 logPow:
[ 14 ] 2,4-Di-tert-butyl-6-(5-chloro-2H-1,2,3-benzotriazol-2-yl)phenol
                                                      Molecular formula:
                                                                              C_{20}H_{24}ClN_3O
                                                                   CAS:
                                                                              3864-99-1
                                                                  ENCS:
                                                                               5-3581, 5-3605
                             HO
                                                                              357.88
                                                                    MW:
                                                                     mp:
                                                                              Uncertain
                                                                      bp:
                                                                              Uncertain
                                                                     SW:
                                                                              Uncertain
                                                         Specific gravity:
                                                                              Uncertain
                                                                 logPow:
                                                                              Uncertain
[ 15 ] N,N-Dimethyldithiocarbamamic acid and its salt
                                                      Molecular formula:
                                                                              C_3H_7NS_2
                                                                    CAS:
                                                                              79-45-8 etc.
                                                                  ENCS:
                                                                               2-1798
                                                                               121.22
                                                                    MW:
                                                                              Uncertain
                                                                     mp:
                                                                      bp:
                                                                              Uncertain
                                                                     SW:
                                                                               Uncertain
                                                         Specific gravity:
                                                                              Uncertain
                                                             logPow:
                                                                              Uncertain
[ 15-1 ] Zinc bis(N,N-dimethyldithiocarbamate) (synonym: Ziram)
                                                      Molecular formula:
                                                                              C_6H_{12}N_2S_4Zn
                                                                    CAS:
                                                                              137-30-4
                                                                  ENCS:
                                                                               2-1833, 9-607
                                                                    MW:
                                                                               305.84
                                                                              249^{26}
                                                                     mp:
                                                                      bp:
                                                                              Uncertain
                                                                     SW:
                                                                              65mg/L (25
                                                                               1.66 (25/4 )<sup>27)</sup>
                                                         Specific gravity:
                                                                             1.23<sup>10)</sup>
                                                              logPow:
[ 15-2 ] N,N'-Ethylenebis(thiocarbamoylthiozinc) bis(N,N-dimethyldithiocarbamate) (synonym: Polycarbamate)
                                                      Molecular formula:
                                                                              C_{10}H_{18}N_4S_8Zn_2
                                                                              64440-88-6
                                                                    CAS:
                                                                  ENCS:
                                                                              2-1848
                                                                    MW:
                                                                              581.61
                                                                              Uncertain
                                                                     mp:
                                                                              Uncertain
                                                                     bp:
                                                                     SW:
                                                                               Uncertain
                                                         Specific gravity:
                                                                              Uncertain
                                                                 logPow:
                                                                              Uncertain
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[ 16 ] N,N-Dimethyldodecylamine N-oxide		
	Molecular formula:	C <sub>14</sub> H <sub>31</sub> NO
	CAS:	1643-20-5
	ENCS:	2-198
O→N—C <sub>12</sub> H <sub>25</sub>	MW:	229.40 132 ~ 133 16)
O→N—C <sub>12</sub> H <sub>25</sub>	mp:	132 133
	bp: SW:	Uncertain
		$190g/L (25)^{17}$
	Specific gravity:	Uncertain 4.67 <sup>4)</sup>
[ 17 ] N,N-Dimethylformamide	logPow:	4.07
[ 17 ] 1v,1v-Dimensynormalinde		
	Molecular formula:	C <sub>3</sub> H <sub>7</sub> NO
	CAS:	68-12-2
	ENCS:	2-680
Ņ	MW:	73.09
	mp:	-61 <sup>5)</sup>
	bp:	153 5)
	SW:	Readily soluble <sup>5)</sup>
<b>н</b> О	Specific gravity:	$0.945(20/4)^{5}$
	logPow:	-1.01 <sup>9)</sup>
[ 18 ] Hydrogenated terphenyl		
	Molecular formula:	$C_{18}H_{(14+i)}$ (i = l+m+n = 1 ~ 14)
$H_1 \stackrel{\checkmark}{\longrightarrow} \longrightarrow \longrightarrow$	CAS:	61788-32-7
	ENCS:	4-41
H <sub>m</sub>	MW:	232.32 ~ 248.45
	mp:	dependent on the molecule
	bp:	dependent on the molecule
H <sub>m</sub>	SW:	dependent on the molecule
$/\!\!/ H_n /\!\!/ H_n$	Specific gravity:	dependent on the molecule
l+m+n =1 ~ 14	logPow:	dependent on the molecule
[ 19 ] O,O-Diethyl O-2-isopropyl-6-methyl-4-pyri	midinyl phosphorothioate	(synonym: Diazinon)
	Molecular formula:	$C_{12}H_{21}N_2O_3PS$
S	CAS:	333-41-5
Ĭ	ENCS:	5-923
O-P-O /	MW:	304.35
$N = \langle  \rangle$	mp:	<25 5)
N=\ O	bp:	$83 \sim 84  (0.002 \text{mmHg})^{5}$
	SW:	40mg/L (25 ) <sup>18)</sup>
N—//	Specific gravity:	$1.12 (20/4)^{5}$
<b>''</b>	logPow:	3.81 <sup>9)</sup>
[ 20 ] O,O-Dimethyl O-3-methyl-4-nitrophenyl ph	osphorothioate (synonym:	Fenitrothion or MFP)
[20] O,O Dimensyl O 3 meanyl 4 introphenyl pin		
S	Molecular formula:	$C_9H_{12}NO_5PS$
	CAS:	122-14-5
0-P-0	ENCS:	3-2616
	MW:	277.23 3.4 <sup>19)</sup>
	mp:	
	bp:	118 $(0.05 \text{mmHg})^{5}$
<u> </u>	SW:	$38 \text{mg/L} (25)^{20}$
O <sub>2</sub> N	Specific gravity:	$1.32 (20/4)^{5}$
<u> </u>	logPow:	3.309)

[ 21 ] Tatrachloroicophthalonitrila (synonym: Chlorothalonil or TDN)	
[ 21 ] Tetrachloroisophthalonitrile (synonym: Chlorothalonil or TPN)	
Molecular formula	$C_8Cl_4N_2$
CAS ENCS	
CI CI MW	
mp	10)
bp	
cw.	
CI CN Specific gravity	$1.8 (25/4)^{10}$
CI logPow	
[ 22 ] Tetrahydrofuran	. 3.03
Molecular formula	$C_4H_8O$
CAS	: 109-99-9
ENCS	
MW	
<b>O</b> mp	
bp	: 65 <sup>6)</sup>
SW	
Specific gravity	
logPow	
[ 23 ] Trichloroacetaldehyde	
[ 25 ] Themstoneounderly at	
Molecular formula	: C <sub>2</sub> HCl <sub>3</sub> O
CI CAS	
ENCS	
MW	
CI———— mp	
bp	
SW O	· /
Specific gravity	
logPow	: 0.99 <sup>9)</sup>
[ 24 ] Trichloronitromethane (synonym: Chloropicrin)	
	· CCL-NO.
Molecular formula	* =
Molecular formula CAS	: 76-06-2
Molecular formula	: 76-06-2 : 2-199
Molecular formula CAS ENCS MW	: 76-06-2 : 2-199 : 164.38
CI Molecular formula CAS ENCS MW mp	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup>
CI Molecular formula CAS ENCS MW mp bp	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup> : 112 <sup>5)</sup>
CI NO <sub>2</sub> Molecular formula CAS ENCS MW mp bp SW	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup> : 112 <sup>5)</sup> : 1,620mg/L (25 ) <sup>7)</sup>
CI NO <sub>2</sub> Molecular formula CAS ENCS MW mp bp SW Specific gravity	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup> : 112 <sup>5)</sup> : 1,620mg/L (25 ) <sup>7)</sup> : 1.66 (20/4 ) <sup>5)</sup>
CI NO <sub>2</sub> Molecular formula CAS ENCS MW mp bp SW Specific gravity logPow	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5</sup> : 112 <sup>5</sup> : 1,620mg/L (25 ) <sup>7</sup> : 1.66 (20/4 ) <sup>5</sup>
CI NO <sub>2</sub> Molecular formula CAS ENCS MW mp bp SW Specific gravity	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup> : 112 <sup>5)</sup> : 1,620mg/L (25 ) <sup>7)</sup> : 1.66 (20/4 ) <sup>5)</sup>
CI NO <sub>2</sub> Molecular formula CAS ENCS MW mp bp SW Specific gravity logPow	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5</sup> : 112 <sup>5</sup> : 1,620mg/L (25 ) <sup>7</sup> : 1.66 (20/4 ) <sup>5</sup> : 2.09 <sup>9</sup> : C <sub>12</sub> H <sub>7</sub> Cl <sub>2</sub> NO <sub>3</sub>
CI CI ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5</sup> : 112 <sup>5</sup> : 1,620mg/L (25 ) <sup>7</sup> : 1.66 (20/4 ) <sup>5</sup> : 2.09 <sup>9</sup> : C <sub>12</sub> H <sub>7</sub> Cl <sub>2</sub> NO <sub>3</sub> : 1836-75-5
CI CI ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS ENCS  Molecular formula CAS ENCS	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5</sup> : 112 <sup>5</sup> : 1,620mg/L (25 ) <sup>7</sup> : 1.66 (20/4 ) <sup>5</sup> : 2.09 <sup>9</sup> : C <sub>12</sub> H <sub>7</sub> Cl <sub>2</sub> NO <sub>3</sub> : 1836-75-5 : 3-658
CI CAS ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS ENCS MW	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup> : 112 <sup>5)</sup> : 1,620mg/L (25 ) <sup>7)</sup> : 1.66 (20/4 ) <sup>5)</sup> : 2.09 <sup>9)</sup> : C <sub>12</sub> H <sub>7</sub> Cl <sub>2</sub> NO <sub>3</sub> : 1836-75-5 : 3-658 : 284.09
CI CAS ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS ENCS MW mp	: 76-06-2 : 2-199 : 164.38 : -64 <sup>5)</sup> : 112 <sup>5)</sup> : 1,620mg/L (25 ) <sup>7)</sup> : 1.66 (20/4 ) <sup>5)</sup> : 2.09 <sup>9)</sup> : C <sub>12</sub> H <sub>7</sub> Cl <sub>2</sub> NO <sub>3</sub> : 1836-75-5 : 3-658 : 284.09 : 70 ~ 71 <sup>23)</sup>
CI CAS ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS ENCS MW mp bp	: $76\text{-}06\text{-}2$ : $2\text{-}199$ : $164.38$ : $-64^{-5)}$ : $112^{-5)}$ : $1,620\text{mg/L} (25^{-})^{7)}$ : $1.66 (20/4^{-})^{5)}$ : $2.09^{9)}$ : $C_{12}H_7Cl_2NO_3$ : $1836\text{-}75\text{-}5$ : $3\text{-}658$ : $284.09$ : $70 \sim 71^{-23)}$ : Uncertain
CI CI ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS ENCS MW mp bp SW Specific gravity logPow  Molecular formula CAS ENCS MW mp bp SW	: $76\text{-}06\text{-}2$ : $2\text{-}199$ : $164.38$ : $-64^{-5}$ : $112^{-5}$ : $1,620\text{mg/L} (25^{-})^{7}$ : $1.66 (20/4^{-})^{5}$ : $2.09^{9}$ : $C_{12}H_7Cl_2NO_3$ : $1836\text{-}75\text{-}5$ : $3\text{-}658$ : $284.09$ : $70 \sim 71^{-23}$ : Uncertain : $1\text{mg/L} (22^{-})^{7}$
CI CAS ENCS MW mp bp SW Specific gravity logPow  [ 25 ] Nitrofen (synonym: NIP)  Molecular formula CAS ENCS MW mp bp hp	: $76\text{-}06\text{-}2$ : $2\text{-}199$ : $164.38$ : $-64^{-5}$ : $112^{-5}$ : $1,620\text{mg/L} (25^{-})^{7}$ : $1.66 (20/4^{-})^{5}$ : $2.09^{9}$ : $C_{12}\text{H}_7\text{Cl}_2\text{NO}_3$ : $1836\text{-}75\text{-}5$ : $3\text{-}658$ : $284.09$ : $70 \sim 71^{-23}$ : Uncertain : $1\text{mg/L} (22^{-})^{7}$ : $1.3^{24}$

[ 26 ] 1,1-Bis( <i>tert</i> -butyldioxy)-3,3,5-trimethylcycloh	nexane	
	Molecular formula:	$C_{17}H_{34}O_4$
	CAS:	6731-36-8
	ENCS:	3-2341
0-0, 0-0-	MW:	302.45
	mp:	-30 <sup>6)</sup>
	bp:	63 6)
	SW:	$0.6 \text{mg/L}^{12)}$
	Specific gravity:	Uncertain
	logPow:	$6.53^{12)}$
[ 27 ] Hydrazine		
	Molecular formula:	$H_4N_2$
	CAS:	302-01-2
$H_2NNH_2$	ENCS:	1-374
11214 14112	MW:	32.05 2.0 <sup>5)</sup>
	mp: bp:	113.5 5)
	SW:	Readily soluble <sup>28)</sup>
	Specific gravity:	1.00 $(25/4)^{5}$
	logPow:	-2.07 <sup>9)</sup>
[ 28 ] 1-Butanol	logi ow.	-2.01
[20] I Bullinoi		
	Molecular formula:	$C_4H_{10}O$
	CAS:	71-36-3
	ENCS:	2-3049
·	MW:	74.12
011	mp:	-89.8 <sup>6)</sup>
· · ·	bp:	117.7 6)
	SW:	63.2g/L (25 ) <sup>29)</sup>
	Specific gravity:	$0.810(20/4)^{6}$
500 J D C J	logPow:	$0.88^{9)}$
[ 29 ] Furfural		
	Molecular formula:	$C_5H_4O_2$
	CAS:	98-01-1
// <b>'</b>	ENCS:	5-40
//	MW:	96.08
H	mp:	-38.1 <sup>6)</sup>
	bp:	161.7 6)
	SW:	$74.1 \text{g/L} (25)^{7}$
	Specific gravity:	$1.16(25/4)^{6}$
0	logPow:	$0.41^{9)}$
[ 30 ] 2-(2-Benzothiazolyloxy)- <i>N</i> -methylacetanilide	(synonym: Mefenacet)	
, , , , , , , , , , , , , , , , ,		
	Molecular formula:	$C_{16}H_{14}N_2O_2S$
l 🐟 .s	CAS:	73250-68-7
	ENCS:	Uncertain
	MW:	298.36
	mp:	134.8 5)
	bp:	Uncertain
Ö	SW:	$4 \text{mg/L} (20)^{20}$
	Specific gravity:	Uncertain 3.23 <sup>10)</sup>
	logPow:	3.23

[ 31 ] 2-(2 <i>H</i> -1,2,3-Benzotriazol-2-yl)-4,6-di- <i>tert</i> -bu	tylphenol	
но	Molecular formula: CAS: ENCS:	C <sub>20</sub> H <sub>25</sub> N <sub>3</sub> O 3846-71-7 5-3580, 5-3604
N N	MW:	323.43
	mp:	Uncertain
N─\ /	bp:	Uncertain
N W	SW:	Uncertain
	Specific gravity: logPow:	Uncertain 6.23 <sup>9)</sup>
	logPow:	0.23
[ 32 ] Methyl methacrylate		
	M 1 1 C 1	CHO
	Molecular formula: CAS:	C <sub>5</sub> H <sub>8</sub> O <sub>2</sub> 80-62-6
) 	ENCS:	2-1036
	MW:	100.12
	mp:	-48 <sup>6)</sup>
) Y	bp:	105.5 6)
	SW:	15.9mg/L (25 ) <sup>7)</sup>
1	Specific gravity:	$0.944 (20/4)^{6}$
	logPow:	1.389)
[ 33 ] 2-(1-Methylethoxy)ethanol		
	Molecular formula:	$C_5H_{12}O_2$
	CAS:	109-59-1
	ENCS:	2-410, 2-2424
\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \	MW:	104.15
У У ОН	mp:	Uncertain
	bp:	145 6)
ı	SW:	Readily soluble <sup>6)</sup>
	Specific gravity:	$0.903 (20/4)^{6}  0.05^{9}$
[ 34 ] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furanyl N	logPow:	
[ 34 ] 2,3-Diffydro-2,2-difficulty1-7-ochzo[o]furaliy17v		yiii. Carooruran)
0	Molecular formula:	$C_{12}H_{15}NO_3$
<b>)</b>	CAS:	1563-66-2
	ENCS:	5-5540 221 25
0 N	MW:	221.25 151 <sup>6)</sup>
Ĭ Ĥ	mp: bp:	Uncertain
0-	SW:	320mg/L (25 ) <sup>18)</sup>
	Specific gravity:	1.18 (20 ) <sup>6)</sup>
	logPow:	2.32 <sup>9)</sup>
[ 25 ] 2 sag Putulphonyl N mathylasakamata (	num: Eanabussek se DDI	AC)
[ 35 ] 2-sec-Butylphenyl N-methylcarbamate; (synon	nym: renoducard or BPN	
•	Molecular formula:	$C_{12}H_{17}NO_2$
)=	CAS:	3766-81-2
	ENCS:	3-2211
O N H	MW:	207.27 32 ~ 33 10)
	mp:	32 - 33
	bp: SW:	00 03
) Y	Specific gravity:	420mg/L (20 ) <sup>10)</sup> Uncertain
	logPow:	2.78 <sup>9)</sup>
	logi ow.	2.70

[ 36 ] α-Methylstyrene	
Molecular formula:	$C_9H_{10}$
CAS:	98-83-9
ENCS:	3-5, 3-8
MW:	118.18
mp:	-23.2 6)
bp:	165.4 6)
SW:	$116 \text{mg/L} (25)^{7}$
Specific gravity:	$0.911 (20/4)^{6}$
logPow:	$3.48^{9)}$
[ 37 ] Dimethyl 2,2-dichlorovinyl phosphate (synonym: Dichlorvos or DDVP)	
Molecular formula:	$C_4H_7Cl_2O_4P$
CAS:	62-73-7
ENCS:	2-3224
MW:	220.98
mp:	<-60 <sup>26)</sup>
CI O—P—O bp:	140 (20mmHg) <sup>31)</sup>
SW:	$8g/L (20)^{2}$
Specific gravity:	$1.42 (25/4)^{6}$
logPow:	$1.47^{32}$
[ 38 ] Tributyl phosphate	
Molecular formula:	$C_{12}H_{27}O_4P$
CAS:	126-73-8
ENCS:	2-2021
MW:	266.31 80 <sup>33)</sup>
mp:	<-80
bp:	289 6)
SW:	280mg/L (25 ) <sup>34)</sup>
Specific gravity:	0.973 <sup>6</sup> )
\ logPow:	$4.00^{9}$

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### 3. Surveyed site and procedure

In the Detailed Environmental Survey (of surface water, sediment, wildlife, and air), the sampling and analysis of specimens was entrusted to prefectural governments and government-designated cities across Japan, and some specimens were analysed by private analytical laboratories. In the survey of surface water for target chemicals [2], [6], and [15] (mainly used as pesticides), the water was sampled taking into consideration the time of pesticide spraying.

In the Detailed Environmental Survey (of food), specimens were sampled and analysed by private analytical laboratories.

- •Survey by market basket method: Market food was surveyed in two regions: Kanto and Kinki. Based on the FY 2004 National Health and Nutrition Survey and its basic data on food intake, according to food groups, food items are classified into 14 groups, and the food contents and their quantities were determined on a daily basis. The food items purchased from grocery stores were treated by the standard method and then rapidly mixed on food group bases and homogenised using a stainless-steel mixer. The mixture is stored in a sealed container, and then used as the sample.
- Survey for domestic food by *Kagezen* method: Japan is divided into 10 regions with 5 selected households per region. The specimens are sampled from meals in three days for each household in a *kagezen* method (the subject household is asked to prepare an extra serving for survey). The specimens from meals in one day (three meals + between-meal snack + beverages) are transferred to a sampling container (acetone-cleaned lidded stainless-steel 3L container), and the closed container is stored in a refrigerator. The meals sampled for 3 consecutive days are treated together, partially thawed, and then rapidly mixed and homogenised using a stainless-steel mixer, taking the greatest care against contamination. The mixture is stored in a sealed container, and then used as the sample.
- Survey for individual food: Instant foods and restaurant food are prepared by the same method as that used for *Kagezen* food.

# (1) Organisations responsible for sampling

Local	Organisations responsible for sempling	Cymfo	Surveye	veyed media		
communities	Organisations responsible for sampling	Surface water	Sediment	Wildlife	Air	
Hokkaido	Hokkaido Institute of Environmental Sciences					
Sapporo City	Sapporo City Institute of Public Health					
Iwate Pref.	Research Institute for Environmental Sciences and Public Health of Iwate Prefecture					
Miyagi Pref.	Miyagi Prefectural Institute of Public Health and Environment					
Sendai City	Sendai City Institute of Public Health					
Yamagata Pref.	Environmental Science Research Center of Yamagata Prefecture					
Ibaraki Pref.	Ibaraki Kasumigaura Environmental Science Center					
Tochigi Pref.	Tochigi Prefectural Institute of Public Health and Environmental Science					
Saitama Pref.	Center for Environmental Science in Saitama					
Chiba Pref.	Chiba Prefectural Environmental Research Center					
Tokyo	Tokyo Metropolitan Research Institute for Environmental Protection					
Kanagawa Pref.	Kanagawa Environmental Research Center					
Yokohama City	Yokohama Environmental Science Research Institute					
Kawasaki City	Kawasaki Municipal Research Institute for Environmental Protection					
Niigata Pref.	Niigata Prefectural Institute of Public Health and Environmental Sciences					
Ishikawa Pref.	Ishikawa Prefectural Institute of Public Health and Environmental Science					
Nagano Pref.	Nagano Environmental Conservation Research Institute					
Shizuoka Pref.	Shizuoka Institute of Environment and Hygiene					
Gifu Pref.	Gifu Prefectural Research Institute for Health and Environmental Sciences					
Aichi Pref.	Aichi Environmental Research Center					
Mie Pref.	Mie Prefectural Science and Technology Promotion Center					
Shiga Pref.	Lake Biwa Environmental Research Institute					
Kyoto Pref.	Kyoto Prefectural Institute of Public Health and Environment					
Kyoto City	Kyoto City Institute of Health and Environmental Sciences					
Osaka Pref.	Osaka Prefecture Environmental Pollution Control Center					
Osaka City	Osaka City Institute of Public Health and Environmental Sciences					
Hyogo Pref.	Hyogo Prefectural Institute of Public Health and Environmental Sciences					
Kobe City	Environmental Conservation and Guidance Division, Environment Bureau					
Wakayama Pref.	Wakayama Prefectural Research Center of Environment and Public Health					
Shimane Pref.	Shimane Prefectural Institute of Public Health and Environmental Science					
Okayama Pref.	Okayama Prefectural Institute for Environmental Science and Public Health					
Hiroshima Pref.	Hiroshima Prefectural Institute of Public Health and Environment					
Hiroshima City	Hiroshima City Institute of Public Health					
Yamaguchi Pref.	Yamaguchi Prefectural Institute of Public Health and Environment					
Tokushima Pref.	Tokushima Prefectural Institute of Public Health and Environmental Sciences					
Kagawa Pref.	Kagawa Prefectural Research Institute for Environmental Sciences and Public Health					

		Survey	ed media	
Organisations responsible for sampling	Surface	Sadiment	Wildlife	Air
	water Sediment Wildlife  te of Public Health and  lth and Environmental  or Hygiene and the  umental Research Center	All		
Ehime Prefectural Institute of Public Health and				
Environmental Science				
Fukuoka Institute of Health and Environmental				
Science				
Fukuoka City Institute for Hygiene and the				
Environment				
Saga Prefectural Environmental Research Center				
Kagoshima Prefectural Institute for Environmental				
Research and Public Health				
	Ehime Prefectural Institute of Public Health and Environmental Science Fukuoka Institute of Health and Environmental Science Fukuoka City Institute for Hygiene and the Environment Saga Prefectural Environmental Research Center Kagoshima Prefectural Institute for Environmental	water  Ehime Prefectural Institute of Public Health and Environmental Science  Fukuoka Institute of Health and Environmental Science  Fukuoka City Institute for Hygiene and the Environment  Saga Prefectural Environmental Research Center  Kagoshima Prefectural Institute for Environmental	Organisations responsible for sampling  Surface water  Ehime Prefectural Institute of Public Health and Environmental Science  Fukuoka Institute of Health and Environmental Science  Fukuoka City Institute for Hygiene and the Environment  Saga Prefectural Environmental Research Center  Kagoshima Prefectural Institute for Environmental	Organisations responsible for sampling  Surface water  Ehime Prefectural Institute of Public Health and Environmental Science  Fukuoka Institute of Health and Environmental Science  Fukuoka City Institute for Hygiene and the Environment  Saga Prefectural Environmental Research Center  Kagoshima Prefectural Institute for Environmental

(Note) Organisations responsible for sampling are described by their official names in FY 2006.

### (2) Surveyed sites (or areas) and target chemicals

Surveyed sites and target chemicals for surface water are shown in Table 2-1-1 and Figure 2-1-1. Surveyed sites and target chemicals for sediment are shown in Table 2-1-2 and Figure 2-1-1. Surveyed areas and target chemicals for wildlife are shown in Table 2-1-3 and Figure 2-1-2. Surveyed sites and target chemicals for air are shown in Table 2-1-4 and Figure 2-1-3. The breakdown is summarized as follows.

Surveyed media	Numbers of local communities	Numbers of target chemicals (groups)	Numbers of surveyed sites (or areas)	Numbers of samples at a surveyed site (or area)
Surface water	29	22	37	3 1
Sediment	10	7	13	3
Wildlife	14	11	15	3
Air	14	7	15	3
Surface water	-	2	-	178 <sup>2</sup>

(Note 1) For target chemicals [2], [6] and [15], 9 specimens were sampled during 3 days taking into consideration the time of pesticide spraying at Katta-bashi Bridge, Riv. Naka (Hitachinaka City) in Ibaraki Pref., Sakae-bashi Bridge, Riv. Tone (Tone Town) in Ibaraki Pref., Kamenoko-bashi Bridge, Riv. Tsurumi (Yokohama City) in Yokohama City, Mouth of Riv. Yamato (Osaka City, Sakai City) in Osaka Pref., and Mishima area, Riv. Iwamatsu (Uwajima City) in Ehime Pref.

(Note 2) Total sample number

Table 2-1-1 (1/2) List of surveyed sites (surface water) and target chemicals in the Detailed Environmental Survey in EV 2006

Local	Surveyed areas	Target chemicals  [1] [2] [3] [4] [5] [6] [7] [10] [11] [12] [1										
communities	•	[1]	[2]	[3]	[4]	[5]	[6]	[7]	[10]	[11]	[12]	[14]
Hokkaido	Ishikarikakokyo Bridge, Mouth of Riv. Ishikari (Ishikari City)											
Sapporo City	Azuma-bashi Bridge, Riv. Toyohira (Sapporo city)											
Iwate Pref.	Riv. Toyosawa (Hanamaki City)											
Miyagi Pref.	Miya-ohashi Bridge, Riv. Matsukawa (Zao Town)											
Sendai City	Hirose-ohashi Bridge, Riv. Hirose (Sendai City)											
Yamagata Pref.	Mouth of Riv. Mogami (Sakata City)											
Ibaraki Pref.	Katta-bashi Bridge, Riv. Naka (Hitachinaka City)											
	Sakae-bashi Bridge, Riv. Tone (Tone Town)											
Tochigi Pref.	Riv. Tagawa (Utsunomiya City)											
Saitama Pref.	Shiki-ohasi Bridge, Riv. Yanase (Shiki City)											
	Asahi-bashi Bridge, Riv. Shingashi (Kawagoe City)											
	Kachi-hashi Bridge, Riv. Ichino (Yoshimi Town)											
Chiba Pref.	Coast of Ichihara and Anegasaki											
Yokohama City	Kamenoko-bashi Bridge, Riv. Tsurumi (Yokohama City)											
Kawasaki City	Mouth of Riv. Tama (Kawasaki City) Keihin Canal in Port of Kawasaki											
Niigata Pref.	Lower Riv. Shinano (Niigata City)											
Ishikawa Pref.	Mouth of Riv. Sai (Kanazawa City)											
Nagano Pref.	Lake Suwa (center)											
Aichi Pref.	Nagoya Port											
Mie Pref.	Yokkaichi Port											
Kyoto City	Miyamae Bridge, Riv. Katsura (Kyoto City)											
Osaka Pref.	Mouth of Riv. Yamato (Sakai City)											
	Mouth of Riv. Yamato (Osaka City, Sakai City)											
Osaka City	Kema-bashi Bridge, Riv. Oh-kawa (Osaka City)											
	Osaka Port											
Hyogo Pref.	Ibogawa-ohashi Bridge, Riv. Ibo (Tatsuno City)											
Kobe City	Kobe Port (center)											
Wakayama Pref.	Kinokawa-ohashi Bridge, Mouth of Riv. Kinokawa (Wakayama City)											
Okayama Pref.	Offshore of Mizushima											
Yamaguchi Pref.	Tokuyama Bay											
	Offshore of Hagi											
Ehime Pref.	Mishima area, Riv. Iwamatsu (Uwajima City)											
Fukuoka Pref.	Kabura-bashi Bridge, River Raizan				-							
i ukuoka i ici.	(Maebaru City)											
	Offshore of Omuta											
Fukuoka City	Hakata Bay											
Saga Pref.	Imari Bay											
20001101.	man Day	l	<u> </u>		<u> </u>		l	<u> </u>				

<sup>[1] 2,2&#</sup>x27;-Azobisisobutyronitrile, [2] 2-Amino-4-[hydroxy(methyl)phosphinoyl]butyric acid (synonym: Glufosinate),

<sup>[3]</sup> m-Aminophenol, [4] Aldicarb, [5] O-Ethyl O-4-nitrophenyl phenylphosphonothioate (synonym: EPN),

<sup>[6]</sup> N,N'-Ethylenebis(dithiocarbamamic acid) and its salt, [7] 2,6-Xylenol, [10] Diisopropylnaphthalene, [11] S-4-Chlorobenzyl N,N-diethylthiocarbamate (synonym: Thiobencarb), [12] 3-(3,4-Dichlorophenyl)-1,1-dimethylurea (synonym: Diuron or DCMU), [14] 2,4-Di-tert-butyl-6-(5-chloro-2H-1,2,3-benzotriazol-2-yl)phenol

Table 2-1-1 (2/2) List of surveyed sites (surface water) and target chemicals in the Detailed Environmental

Local	Surveyed sites						et chen		1	1	1	
communities	•	[15]	[19]	[20]	[21]	[23]	[30]	[31]	[32]	[35]	[37]	[38]
Hokkaido	Ishikarikakokyo Bridge, Mouth of Riv. Ishikari (Ishikari City)											
Sapporo City	Azuma-bashi Bridge, Riv. Toyohira (Sapporo city)											
Iwate Pref.	Riv. Toyosawa (Hanamaki City)											
Miyagi Pref.	Miya-ohashi Bridge, Riv. Matsukawa (Zao Town)											
Sendai City	Hirose-ohashi Bridge, Riv. Hirose (Sendai City)											
Yamagata Pref.	Mouth of Riv. Mogami (Sakata City)											
Ibaraki Pref.	Katta-bashi Bridge, Riv. Naka (Hitachinaka City)											
	Sakae-bashi Bridge, Riv. Tone (Tone Town)											
Tochigi Pref.	Riv. Tagawa (Utsunomiya City)											
Saitama Pref.	Shiki-ohasi Bridge, Riv. Yanase (Shiki City)											
	Asahi-bashi Bridge, Riv. Shingashi (Kawagoe City)											
	Kachi-hashi Bridge, Riv. Ichino (Yoshimi Town)											
Chiba Pref.	Coast of Ichihara and Anegasaki											
Yokohama City	Kamenoko-bashi Bridge, Riv. Tsurumi (Yokohama City)											
Kawasaki City	Mouth of Riv. Tama (Kawasaki City)											
-	Keihin Canal, Port of Kawasaki											
Niigata Pref.	Lower Riv. Shinano (Niigata City)											
Ishikawa Pref.	Mouth of Riv. Sai (Kanazawa City)											
Nagano Pref.	Lake Suwa (center)											
Aichi Pref.	Nagoya Port											
Mie Pref.	Yokkaichi Port											
Kyoto City	Miyamae Bridge, Riv. Katsura (Kyoto City)											
Osaka Pref.	Mouth of Riv. Yamato (Sakai City)											
	Mouth of Riv. Yamato (Osaka City, Sakai City)											
Osaka City	Kema-bashi Bridge, Riv. Oh-kawa (Osaka City)											
	Osaka Port											
Hyogo Pref.	Sakae-bashi Bridge, Riv. Tone (Tone Town)											
Kobe City	Kobe Port (center)											
Wakayama Pref.	Kinokawa-ohashi Bridge, Mouth of Riv. Kinokawa (Wakayama City)											
Okayama Pref.	Offshore of Mizushima											
Yamaguchi Pref.												
	Offshore of Hagi											
Ehime Pref.	Mishima area, Riv. Iwamatsu (Uwajima											
Fukuoka Pref.	City) Kabura-bashi Bridge, River Raizan											
гикиока Ртег.	(Maebaru City)											
	Offshore of Omuta											
Fukuoka City	Hakata Bay											
Saga Pref.	Imari Bay											
	yldithiocarbamamic acid and its salt [10] () ()	L	L	Ļ			<u> </u>	L		<u> </u>	L	

<sup>[15]</sup> *N,N*-Dimethyldithiocarbamamic acid and its salt, [19] *O,O*-Diethyl *O*-2-isopropyl-6-methyl-4-pyrimidinyl phosphorothioate (synonym: Diazinon), [20] *O,O*-Dimethyl *O*-3-methyl-4-nitrophenyl phosphorothioate (synonym: Fenitrothion or MEP),

<sup>[21]</sup> Tetrachloroisophthalonitrile (synonym: Chlorothalonil or TPN), [23] Trichloroacetaldehyde,

<sup>[30] 2-(2-</sup>Benzothiazolyloxy)-N-methylacetanilide (synonym: Mefenacet),

<sup>[31] 2-(2</sup>*H*-1,2,3-Benzotriazol-2-yl)-4,6-di-*tert*-butylphenol, [32]Methyl methacrylate, [35] 2-*sec*-Butylphenyl

N-methylcarbamate; (synonym: Fenobucarb or BPMC), [37]Dimethyl 2,2-dichlorovinyl phosphate (synonym: Dichlorvos or DDVP), [38]Tributyl phosphate

Table 2-1-2 List of surveyed sites (sediment) and target chemicals in the Detailed Environmental Survey in FY 2006

Local	Surveyed sites			Tar	get chemi	cals		
communities	Surveyed sites	[4]	[8]	[14]	[16]	[17]	[31]	[36]
Yamagata Pref.	Mouth of Riv. Mogami (Sakata City)							
Kawasaki City	Mouth of Riv. Tama (Kawasaki City)							
	Keihin Canal in Kawasaki Port							
Shizuoka Pref.	Shimizu Port							
Aichi Pref.	Nagoya Port							
Mie Pref.	Yokkaichi Port							
Osaka Pref.	Mouth of Riv. Yamato (Sakai City)							
Osaka City	Kema-bashi Bridge, Riv. Oh-kawa (Osaka City)							
	Osaka Port							
Hyogo Pref.	Tatsumi-bashi Bridge, Riv. Samondo (Amagasaki City)							
Hiroshima Pref.	Kure Port							
Yamaguchi Pref.	Tokuyama Bay							
	Offshore of Hagi							

<sup>[4]</sup> Aldicarb, [8] Chlorobenzene, [14] 2,4-Di-*tert*-butyl-6-(5-chloro-2*H*-1,2,3-benzotriazol-2-yl)phenol, [16] *N*,*N*-Dimethyldodecylamine *N*-oxide, [17] *N*,*N*-Dimethylformamide,

<sup>[31] 2-(2</sup>H-1,2,3-Benzotriazol-2-yl)-4,6-di-tert-butylphenol, [36] $\alpha$ -Methylstyrene

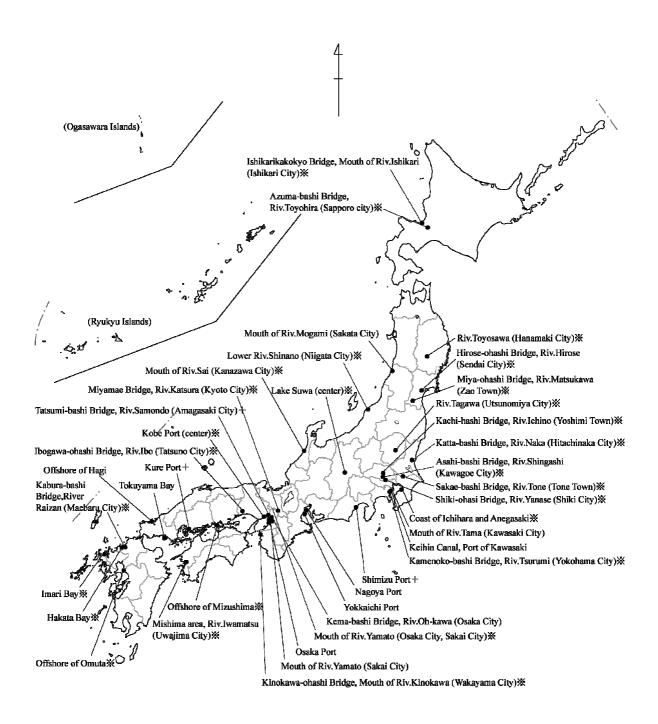


Figure 2-1-1 Surveyed sites (surface water and sediment) in the Detailed Environmental Survey in FY 2006 (Note) means "surface water only." † means "sediment only."

Table 2-1-3 List of surveyed areas (wildlife) and target chemicals in the Detailed Environmental Survey in FY 2006

2006 Local	Cumposed ong	Wildlife angeis -					Targe	et cher	nicals				
communities	Surveyed areas	Wildlife species	[4]	[6]	[12]	[14]	[15]	[18]	[24]	[26]	[27]	[30]	[31]
Hokkaido	Offshore of Japan Sea (offshore of Iwanai)	Greenling (Hexagrammos lagocephalus)											
Iwate Pref.	Yamada Bay	Blue mussel (Mytilus galloprovincialis)											
Miyagi Pref.	Sendai Bay (Matsushima Bay)	Sea bass ( <i>Lateolabrax japonicus</i> )											
Tokyo Met.	Tokyo Bay	Sea bass (Lateolabrax japonicus)											
Yokohama City	Riv. Tsurumi	Carp (Cyprinus carpio)											
Kawasaki City	Offshore of Ogishima Island, Port of Kawasaki	Sea bass (Lateolabrax japonicus)											
Niigata Pref.	Lower Riv. Shinano (Niigata City)	Carp (Cyprinus carpio) and Barbel steed (Hemibarbus barbus)											
Shiga Pref.	Lake Biwa, Riv. Azumi (Takashima City)	Dace (Tribolodon hakonensis)											
Osaka Pref.	Osaka Bay	Sea bass ( <i>Lateolabrax japonicus</i> )											
Shimane Pref.	Shichirui Bay, Shimane Peninsula	Blue mussel (Mytilus galloprovincialis)											
Hiroshima City	Hiroshima Bay	Sea bass ( <i>Lateolabrax japonicus</i> )											
Yamaguchi Pref.	Tokuyama Bay	Striped mullet (Mugil cephalus)											
	Offshore of Hagi	Striped mullet (Mugil cephalus)											
Tokushima Pref.	Naruto	Hard-shelled mussel (Mytilus coruscus)											
Kagoshima Pref.	West Coast of Satsuma Peninsula	Sea bass (Lateolabrax japonicus)			(2.4.D			1) 1					

[4] Aldicarb, [6] *N,N'*-Ethylenebis(dithiocarbamamic acid) and its salt, [12] 3-(3,4-Dichlorophenyl)-1,1-dimethylurea (synonym: Diuron or DCMU), [14] 2,4-Di-*tert*-butyl-6-(5-chloro-2*H*-1,2,3-benzotriazol-2-yl)phenol, [15] *N,N*-Dimethyldithiocarbamamic acid and its salt, [18] Hydrogenated terphenyl, [24] Trichloronitromethane (synonym: Chloropicrin),

[26] 1,1-Bis(*tert*-butyldioxy)-3,3,5-trimethylcyclohexane, [27] Hydrazine, [30] 2-(2-Benzothiazolyloxy)-*N*-methylacetanilide (synonym: Mefenacet), [31] 2-(2*H*-1,2,3-Benzotriazol-2-yl)-4,6-di-*tert*-butylphenol

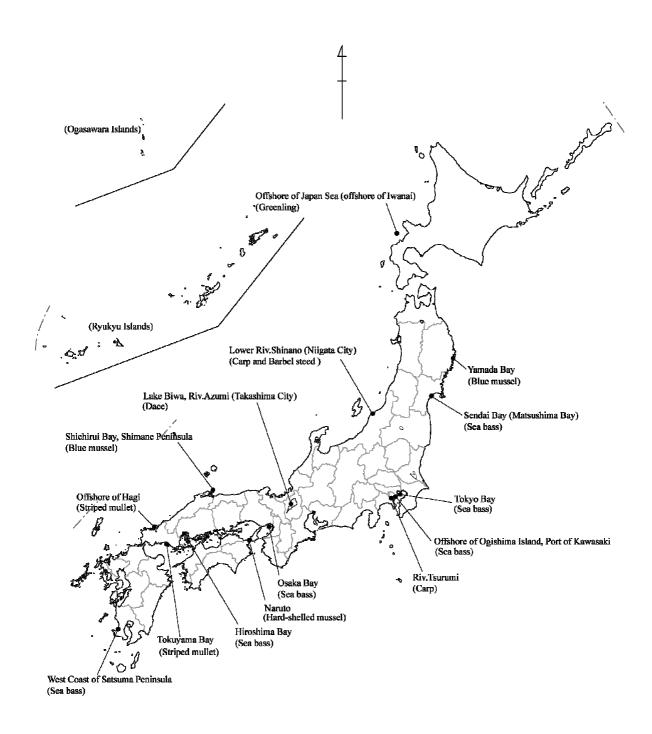


Figure 2-1-2 Surveyed areas (wildlife) in the Detailed Environmental Survey in FY 2006

Table 2-1-4 List of surveyed sites (air) and target chemicals in the Detailed Environmental Survey in FY 2006

Table 2-1-4 List	of surveyed sites (air) and target chemicals in the I	<b>Detaile</b>	d Envi				n FY 2	006
Local	Surveyed sites			Targ	et chem	nicals		
communities	Surveyed sites	[9]	[13]	[22]	[25]	[28]	[29]	[33]
Sapporo City	Sapporo City Institute of Public Health (Sapporo City)							
Saitama Pref.	Center for Environmental Science in Saitama (Kisai							
	Town)							
Chiba Pref.	Ichihara-Matsuzaki Air Quality Monitoring Station							
	(Ichihara City)							
Tokyo Met.	Tokyo Metropolitan Research Institute for							
	Environmental Protection (Koto Ward)							
	Chichijima Island							
Kanagawa Pref.	Kanagawa Environmental Research Center (Hiratsuka							
	City)							
Gifu Pref.	Gifu Prefectural Research Institute for Health and							
	Environmental Sciences (Kakamigahara City)							
Mie Pref.	Mie Prefectural Science and Technology Promotion							
	Center (Yokkaichi City)							
Kyoto Pref.	Kyoto Prefectural Joyo High School (Joyo City)							
Kyoto City	Kyoto City Hall (Kyoto City)							
Osaka Pref.	Osaka Prefecture Environmental Pollution Control							
	Center (Osaka City)							
Hyogo Pref.	Hyogo Prefectural Institute of Public Health and							
	Environmental Sciences (Kobe City)							
Wakayama Pref.	Wakayama Prefectural Research Center of Environment							
	and Public Health (Wakayama City)							
Yamaguchi Pref.	Yamaguchi Prefectural Institute of Public Health and							
	Environment (Yamaguchi City)							
Kagawa Pref.	Takamatsu Joint Prefectural Government Building							
	(Takamatsu City)							

[9] Isobutyl acetate, [13] 2,6-Dichlorobenzonitrile (synonym: Dichlobenil), [22] Tetrahydrofuran, [25] Nitrofen (synonym: NIP), [28] 1-Butanol, [29] Furfural, [33] 2-(1-Methylethoxy)ethanol



Figure 2-1-3 Surveyed sites (air) in the Detailed Environmental Survey in FY 2006

 $Table \ 2\text{-}1\text{-}5 \ List \ of surveyed \ regions \ (food) \ and \ target \ chemicals \ in \ the \ Detailed \ Environmental \ Survey \ in \ FY \ 2006$ 

Survey by market basket method

	For market basket	ood			Targ	get Chemicals
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furany 1 <i>N</i> -methylcarbamate (synonym: Carbofuran)
	Rice grains	Well-milled rice (1)	Kanto	163.1	_	
			Kinki	159.3	_	
Group		Well-milled rice (2)	Kanto	163.1	one specimen from	one specimen from one
1			Kinki	159.3	one mixed sample	mixed sample per survey
	Processed rice	Rice cake	Kanto	5.2	per survey region	region
			Kinki	5.1		
	Wheat flours	Soft or hard wheat flours	Kanto	8.1	4	
	D 1 ( )	3371.4 4 11 1 1	Kinki	8.5	_	
	Breads (except sweet buns)	White table bread	Kanto	17.2	-	
	sweet buils)	E 11 1	Kinki	18.7	-	
		French bread	Kanto	17.2	4	
	C	D £11. 1:41	Kinki	18.7	4	
	Sweet buns	Bun filled with sweat bean paste	Kanto	6.6	_	
	III CI:	*	Kinki	7.2	_	
	Udon, Chinese noodles	Raw udon	Kanto	12.3	_	
	lioodies	D CI: II	Kinki	13.3	-	
		Raw Chinese noodles	Kanto	12.2	-	
		D: 1	Kinki	13.3	_	
		Dried somen	Kanto	12.2	_	
		7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7 7	Kinki	13.3	4	
	Instant Chinese	Instant noodle (deep-fried)	Kanto	4.2	_	
	noodles		Kinki	4.6	4	
	Pastas	Macaroni and Spaghetti	Kanto	10.1		
			Kinki	11.0		
	Other wheat	Chio tzu pastry	Kanto	5.2	_	
	processed foods	D 1 1 1 1 1	Kinki	4.4	4	
Cassa	Buckwheat	Raw buckwheat noodles	Kanto	7.1		
Group 2	noodles and processed foods		Kinki	7.7	one specimen from one mixed sample	one specimen from one mixed sample per survey
	Corns and	Corn flakes	Kanto	0.5	per survey region	region
	processed foods		Kinki	0.6		
	Other grains		Kanto	1.2		
			Kinki	2.4		
	Seeds	Sesame seeds	Kanto	1.2		
			Kinki	1.1		
		Oil-roasted and salted	Kanto	1.2		
		peanuts	Kinki	1.1		
	Sweet potatoes	Sweet potatoes	Kanto	7.4		
	and processed foods		Kinki	6.7		
	Potatoes and	Potatoes	Kanto	27.7		
	processed foods		Kinki	26.3		
	Other potatoes	Satoimo (Taro)	Kanto	7.5		
	and processed		Kinki	6.6		
	foods	Japanese yam or Chinese	Kanto	7.6		
	15005	yam	Kinki	6.7		
		Stringy konnyaku	Kanto	6.3		
		(Taro-starch vermicelli)	Kinki	6.9	7	
		1				
	Starches and processed foods	bean-starch vermicelli	Kanto	1.6		

	Fo	od			Targ	et Chemicals	
			Surveyed	Daily intake		[34] 2,3-Dihydro-2,2-	
Groups	Sub-groups	Items	regions	weight	[27] Hydrazine	dimethyl-7-benzo[b]furany	
Croups	Sue groups	1001115		(g)	[27] 11) utuziii	1 <i>N</i> -methylcarbamate	
	Sugars and	Caster sugar	Kanto	6.8		(synonym: Carbofuran)	
	sweeteners	Castel sugai	Kinki	7.6			
	Japanese-style	Sweet azuki-bean jelly	Kanto	5.6			
	confectioneries	5 weet azaki-bean jeny	Kinki	5.5			
					_		
			Kanto Kinki	5.5 5.4	_		
	Cakes and	Shortcakes					
	pastries	Shortcakes	Kanto Kinki	3.9	_		
Group	pastrics	Custard cream puffs	Kanto	3.8			
Group 3		Custaru cream puns	Kinki	3.7	one specimen from one mixed sample	one specimen from one mixed sample per survey	
3	Biscuits	Hard or soft biscuits	Kanto	1.9	per survey region	region	
	Discuits	riald of soft discuits	Kinki	1.9			
	Candies	Fruit juice tasted tablets		0.3			
	Calidies	Truit juice tasted tablets	Kanto Kinki	0.3			
	Other confections	Milk chocolates					
	Other confections	Wilk chocolates	Kanto Kinki	2.8			
		Crisps	Kanto	2.7			
		Crisps	Kinki	2.7	-		
	Butter	Salted butter	Kanto	1.2			
	Dutter	Salled buller	Kinki	1.1			
	Margarine	Soft type margarine	Kanto	1.3			
	Wargarnie	Soft type margarine	Kinki	1.3			
	Vegetable oil	Blended oil	Kanto	3.0	one enceimen from		
Cassan	vegetable on	Dielided Oil	Kinki	2.8			
Group 4		Olive oil	Kiliki	3.0	one specimen from one mixed sample	one specimen from one mixed sample per survey	
		Olive on	Kinki	2.8	per survey region	region	
		Sesame oil	Kanto	3.0			
		Sesame on	Kinki	2.7			
	Animal oil and fat	Lard	Kanto	0.1			
	Anniai on and rat	Laid	Kinki	0.1			
	Soybeans (whole	Boiled soybeans	Kanto	2.1			
	grain) and	Boned soybeans					
	processed foods		Kinki	2.0			
	Bean curds	Hard or soft bean curds	Kanto	35.9			
			Kinki	34.2			
	Deep-fried bean	Deep-fried bean curds	Kanto	4.3			
	curds		Kinki	3.9			
Group		Ganmodoki (Deep-fried	Kanto	3.6	one specimen from	one specimen from one	
5		fritters of mashed bean	Kinki	4.6	one mixed sample	mixed sample per survey	
	N (C) 1	curd and vegetables)			per survey region	region	
	Natto (Steamed and fermented	Sticky Natto	Kanto	6.4	4		
	soybeans)		Kinki	6.1			
	Other soy	Soy milk	Kanto	6.9			
	products	•	Kinki	6.6			
	Other beans and	Boiled red kidney beans	Kanto	1.7			
	processed foods		Kinki	1.6			

	Fo	ood			Targ	get Chemicals		
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furany 1 <i>N</i> -methylcarbamate (synonym: Carbofuran)		
	Strawberries	Strawberries	Kanto	0.2				
			Kinki	0.2				
	Citrus fruits	Oranges	Kanto	16.7				
			Kinki	14.7				
		Grapefruits	Kanto	16.6				
			Kinki	14.6				
	Bananas	Bananas	Kanto	15.2				
			Kinki	13.4				
	Apples	Apples	Kanto 25.7  Kinki 22.6					
			Kinki	22.6				
Group	Other fresh fruits	Watermelons	Kanto	13.0	one specimen from	one specimen from one		
6			Kinki	11.4	one mixed sample	mixed sample per survey		
		Melons	Kanto	13.0	per survey region	region		
			Kinki	11.4				
		Pineapples	Kanto	12.9				
			Kinki	11.3				
	Fruit juice drink	Strawberry jam	Kanto	1.2				
			Kinki	1.1				
		Orange juice	Kanto	8.0				
			Kinki	7.0				
		Apple juice	Kanto	8.0				
			Kinki	7.0				
	Tomatoes	Tomatoes	Kanto	14.2				
			Kinki	12.9				
	Carrots	Carrots	Kanto	17.9				
			Kinki	19.3				
	Spinach	Spinach	Kanto	13.5				
			Kinki	13.7				
	Sweet peppers	Green bell peppers	Kanto	3.3				
Group			Kinki	2.9	one specimen from	one specimen from one		
7	Other green and	Broccoli	Kanto	15.3	one mixed sample	mixed sample per survey		
	yellow vegetables		Kinki	12.8	per survey region	region		
		Asparagus	Kanto	11.9				
			Kinki	10.6				
		Pumpkins	Kanto	11.9				
			Kinki	12.7				
	Vegetable juice	(Carton vegetable juice	Kanto	9.3				
		was used.)	Kinki	7.4				

	Fo	ood			Targ	et Chemicals
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furany l <i>N</i> -methylcarbamate (synonym: Carbofuran)
	Cabbage	Cabbage	Kanto	20.0		
			Kinki	18.6		
	Cucumbers	Cucumbers	Kanto	9.6		
	D .1	D. II	Kinki	8.9		
	Daikon	Daikon	Kanto	27.2		
	0 :	0 :	Kinki	30.4		
	Onions	Onions	Kanto	25.9		
	CI.: 11	CI: 11	Kinki	26.6		
	Chinese cabbage	Chinese cabbage	Kanto	11.7		
	04 1:14	D 1 1 ( 7 1 1)	Kinki	10.0	_	
	Other light coloured	Bamboo-shoots (boiled)	Kanto	8.3		
	vegetables	Doon ammounts	Kinki	7.7		
	regemeres	Bean sprouts	Kanto	7.3		
		Eggplants	Kinki	6.8	_	
		Eggpiants	Kanto Kinki	5.6 5.8		
		Burdocks	Kanto	7.3		
		Durdocks	Kinki	7.6		
		Lettuce	Kanto	8.2		
		Lettuce	Kinki	7.6		
Group	Pickled vegetable leaves	Seasoned Nozawana pickles	Kanto	3.2	one specimen from	one specimen from one
8			Kinki	2.5	one mixed sample	mixed sample per survey
		Salted <i>Chinese</i> cabbage pickles	Kanto	3.1	per survey region	region
			Kinki	2.5		
	Takuan and other	Takuan pickles	Kanto	6.1		
	pickles		Kinki	4.8		
		Fukujin pickles	Kanto	6.0		
		_ ~ ~ .	Kinki	4.8		
	Mushrooms	Raw Shiitake mushroom	Kanto	2.8	_	
		T. In I	Kinki	3.3	_	
		Enokitake mushroom	Kanto	4.6		
		**	Kinki	4.5		
		Hatakeshimeji,	Kanto	4.8		
	G 1	Bunashimeji, and Honshimeji mushroom	Kinki	4.3		
	Seaweeds	Laver	Kanto	3.0	_	
		Voubu (dried)	Kinki	3.1	_	
		Konbu (dried)	Kanto	11.4	_	
		Wakame (dried)	Kinki	11.5	_	
		vv akame (ariea)	Kanto	25.3	_	
		Dried <i>Hijiki</i>	Kinki	25.5	_	
		Difed Hijiki	Kanto	3.6	-	
			Kinki	5.9		

	Fo	ood			Targ	get Chemicals
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2- dimethyl-7-benzo[b]furany 1 <i>N</i> -methylcarbamate (synonym: Carbofuran)
	Japanese Sake	Refined sake	Kanto	11.2		
			Kinki	12.8		
	Beer		Kanto	60.5		
			Kinki	69.1		
	Western liquor	White wine or red wine	Kanto	25.9		
			Kinki	29.6		
	Tea	Green tea (infusion)	Kanto	151.9		
Group			Kinki	173.5	one specimen from	one specimen from one
9		Black tea (infusion)	Kanto	151.9	one mixed sample	mixed sample per survey
			Kinki	173.4	per survey region	region
	Coffee or Cocoa	Coffee	Kanto	121.7		
			Kinki	139.0		
	Other beverages	Coke	Kanto	43.5		
			Kinki	49.7		
			Kanto	43.5		
			Kinki	49.7		

	Fo	ood			Taro	et Chemicals			
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furany 1 <i>N</i> -methylcarbamate (synonym: Carbofuran)			
	Japanese horse	Japanese horse mackerels	Kanto	2.8					
	mackerels and		Kinki	3.2					
	sardines	Japanese pilchard	Kanto	2.8					
			Kinki	3.1					
			Kanto	2.4					
			Kinki	3.0					
		Pacific Saury	Kanto	2.5					
			Kinki	2.9					
	Salmons and		Kanto	3.5					
	trouts		Kinki	3.9					
	Sea breams and		Kanto	3.4					
	flounders		Kinki	3.3					
		Pacific Cod (fillet)	Kanto	3.2	1				
			Kinki	2.7	1				
	Tunas and	Bigeye tuna	Kanto	6.0	one specimen from	one specimen from one			
	marlins	8.7.	Kinki	6.8	one mixed sample	mixed sample per survey			
	Other raw fishes	Yellowtail	Kanto	4.7	per survey region	region			
	Cuiei iu vi iisiies	10110 // 4411	Kinki	5.2					
		Eel (broiled)	Kanto	4.8	-				
		Zer (erenee)	Kinki	5.4					
	Bivalves	Short-necked clams	Kanto	2.0	_				
	Divarves	Short necked claims	Kinki	2.2	_				
		Japanese scallop	Kanto	1.9	†				
		sapanese seanop	Kinki	2.2	†				
	cuttlefishes and octopuses	Japanese common squid	Kanto	2.8	†				
		supunese common squid	Kinki	3.2	-				
Group		Common octopus (boiled)	Kanto	2.8					
10		Common octopus (bonea)	Kinki	3.1					
	Shrimps and		Kanto	5.4					
	crabs		Kinki	6.7	-				
	Seafood (salted,	Salted salmon	Kanto	5.1					
	half dried, and	Salica salilion	Kinki	5.4	+				
	dried)	Salted cod roe	Kinki	4.8	_				
	·	Sanca coa roc	Kinki	5.7	+				
		Japanese horse mackerels (dried fillet)	Kanto	4.9	-				
		(	Kinki	5.5	1				
	Seafood (canned	Canned boiled mackerel	Kanto	1.1	1				
	food)		Kinki	1.3	1				
	·	Canned seasoned tuna	Kanto	1.1					
			Kinki	1.2					
	Seafood	Japanese sand lance	Kanto	0.4	one specimen from	one specimen from one			
	(Tsukudani)	(Tsukudani)	Kinki	0.4	one mixed sample per survey region	mixed sample per survey region			
	Seafood	Steamed Kamaboko	Kanto	2.3	_ relicantely region	1051011			
	(kneaded)		Kinki	2.6	1				
		Baked Chikuwa	Kanto	2.3	1				
			Kinki	2.6	1				
		Натреп	Kanto	2.0	1				
			Kinki	2.4	1				
		Satsumaage	Kanto	2.2	†				
			Kinki	2.5	1				
	Fish hams and	Fish sausages	Kanto	0.4	1				
	sausages	1 1011 00000000	Kinki	0.4	1				
<u> </u>		1	IXIIIKI	U. <del>4</del>					

	Fo	ood			Targ	et Chemicals		
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furany 1 <i>N</i> -methylcarbamate (synonym: Carbofuran)		
	Beef	Brisket	Kanto	6.3				
			Kinki	6.4				
		Ground beef	Kanto	7.0	=			
			Kinki	6.2				
	Pork	Pork thigh	Kanto	13.7	_			
			Kinki	11.7				
		Pork loin	Kanto	14.1				
			Kinki	13.7				
	Hams and	Roast ham	Kanto	4.2				
	sausages		Kinki	3.9				
		Bacon	Kanto	3.7				
Group			Kinki	3.2	one specimen from	one specimen from one		
11		Vienna sausage	Kanto	3.9	one mixed sample	mixed sample per survey		
			Kinki	3.4	per survey region	region		
	Other animal	Lamb shoulder	Kanto	0.3				
	meats		Kinki	0.3				
	Chicken	Chicken thigh	Kanto	17.9				
			Kinki	16.6				
	Other poultries	Canard	Kanto	0.1				
			Kinki	0.1				
	Meets (guts)	Beef liver or pork liver	Kanto	1.2	-			
	,	•	Kinki	1.0	-			
	Eggs	Hen's eggs	Kanto	31.4	1			
	-88*	1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	Kinki	38.3	1			
	Milk	Ordinary liquid milk	Kanto	110.4				
			Kinki	98.1	1			
	Cheese	Processed cheese	Kanto	2.5				
			Kinki	2.2	1			
	Fermented milk	Plain yogurt or sweetened	Kanto	12.6				
	and lactic acid	yogurt	Kinki	11.1				
Group	bacteria beverage	Lactic acid bacteria	Kanto	12.5	one specimen from	one specimen from one		
12		beverage and dairy products	Kinki	11.1	one mixed sample per survey region	mixed sample per survey region		
	Other dairy	Whipping cream	Kanto	4.5	1			
	products	11 0	Kinki	4.0				
		Ice cream	Kanto	4.4	1			
		ice cream		3.9	1			
	i		Kinki	L 2./	1			

	Fo	ood			Targ	get Chemicals
Groups	Sub-groups	Items	Surveyed regions	Daily intake weight (g)	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furany 1 <i>N</i> -methylcarbamate (synonym: Carbofuran)
	Sauce	Worcester sauce	Kanto	2.1		
			Kinki	2.4		
	Soy sauce	Koikuchi	Kanto	16.3		
			Kinki	18.7		
	Salt	Common salt	Kanto	1.5		
			Kinki	1.7		
	Mayonnaise	Mayonnaise	Kanto	3.3		
			Kinki	3.7		
	Miso	Light yellow type	Kanto	11.4		
			Kinki	13.2		
	Other seasonings	Grain vinegar or rice vinegar	Kanto	6.3		
			Kinki	7.2		
		Dressing	Kanto	6.2		
			Kinki	7.2		
Group		Tomato ketchup	Kanto	6.2	one specimen from	one specimen from one
13			Kinki	7.2	one mixed sample	mixed sample per survey
		Mentsuyu	Kanto	6.2	per survey region	region
			Kinki	7.2	]	
			Kanto	6.2		
			Kinki	7.2		
			Kanto	6.2		
			Kinki	7.1		
		Consommé cubes	Kanto	6.2		
			Kinki	7.1		
		Curry roux	Kanto	10.9		
			Kinki	11.5		
		Hash roux	Kanto	10.9		
			Kinki	11.0		
	spices and others	Mustard paste	Kanto	0.1		
		TT 1.	Kinki	0.1	_	
		Wasabi paste	Kanto	0.1		
Group	Drinking water	Tap water	Kinki Kanto	0.1 250	one specimen from	one specimen from one
14			Kinki	250	one mixed sample per survey region	mixed sample per survey region

Survey of nationwide food by Kagezen method

			Target chemicals							
Surveyed areas	Surveyed sites	Numbers of samples	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furanyl N-methylcarbamate (synonym: Carbofuran)						
Hokkaido	Ishikari City	1								
	Sapporo City	1								
	Eniwa City	1								
	Sapporo City	1								
	Otaru City	1								
Miyagi	Sendai City	1								
Pref.	Sendai City	1								
	Sendai City	1								
	Sendai City	1								
	Tagajo City	1								
Tokyo	Hachioji City	1								
Met.	Tama City	1								
	Nerima Ward	1								
	Suginami Ward	1								
	Tama City	1								
Ishikawa	Hakusan City	1								
Pref.	Kanazawa City	1								
	Kanazawa City	1								
	Kanazawa City	1								
	Kahoku City	1								
Nagano Pref.	Ueda City	1								
	Matsumoto City	1								
	Shimoina District	1								
	Saku City	1								
	Nagano City	1								
Aichi	Nishio City	1								
Pref.	Nagoya City	1								
	Kasugai City	1								
	Nagoya City	1								
	Nagoya City	1								
Osaka	Minoh City	1								
Pref.	Izumisano City	1								
	Takaishi City	1								
	Settsu City	1								
	Katano City	1								
Ehime	Matsuyama City	1								
Pref.	Matsuyama City	1								
11011	Matsuyama City	1								
	Matsuyama City	1								
	Matsuyama City	1								
Fukuoka	Fukuoka City	1								
Pref.	Fukuoka City	1		+						
1 101.	Fukuoka City	1								
	Fukuoka City  Fukuoka City	1								
	Kitakyushu City	1								
Olring										
Okinawa Pref.	Urasoe City	1								
riel.	Nakagami District	1								
	Itoman City	1								
	Naha City	1								
	Okinawa City	1								

Survey of individual instant foods

Survey of individual in	istant foods	1	То	rget Chemicals
			Tai	
Frozen pilaf	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furanyl <i>N</i> -methylcarbamate (synonym: Carbofuran)		
	Frozen nilaf	1		(synonym: Carboraran)
	` ′			
Frozen food				
Retort-pouched food				
r				
	1 1 5	_		
		1		
	-	1		
	1			
		1		
		1		
	Instant cup fried noodles	1		
		1		
Instant food	*	1		
		1		
	7	1		
		1		
	Instant bean paste soup	1		
		1		
	Canned seafood (tuna)	1		
		1		
G 16 1		1		
Canned food		1		
	Canned cooked food	1		
75 1.0 . 1				
Bottled food				
	C			
	-	1		
	Preparedfood-4	1		
	Preparedfood-5	1		
Prepared food	Preparedfood-6	1		
	Preparedfood-7	1		
	Preparedfood-8	1		
	Preparedfood-9	1		
	Preparedfood-10	1		
	11cparcu100u-10	1		1

Survey of individual restaurant food

			Target Chemicals					
Types	Contents	Numbers of samples	[27] Hydrazine	[34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furanyl N-methylcarbamate (synonym: Carbofuran)				
Fast food	Hamburger set, etc.	5						
Japanese rice bowl dishes	Rice bowl dishes, etc.	5						
Family restaurant	Set menus	6						
Steak, grilled meat	Set menus	4						
Sushi	Sushi	5						
Ramen, soba, udon, pasta	Noodles	6						
Chinese food	Set menus	4						
Other restaurants	Set menus	5						
Box lunch shop	Box lunch	7						
Bakery	Bread	3						

#### (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 equipments. 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 ).

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 ).

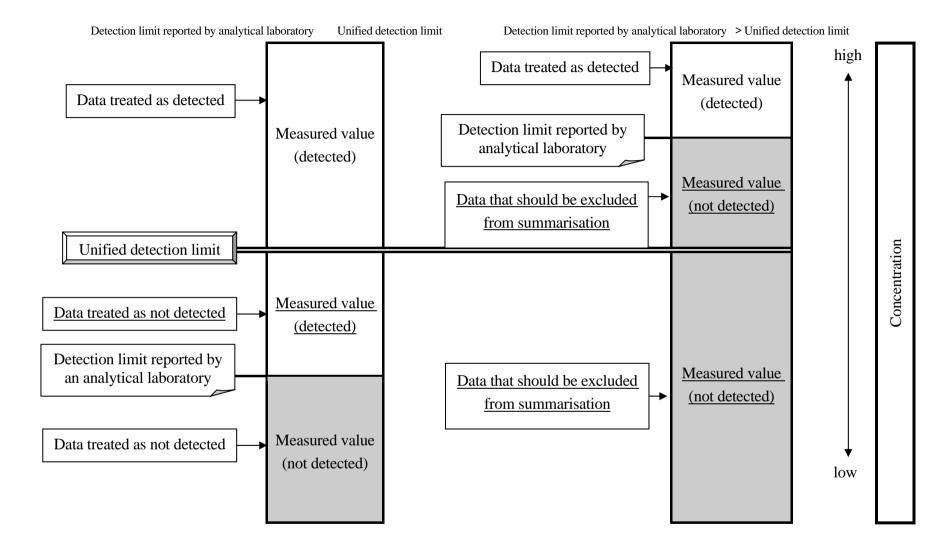
When the instrument detection limit (IDL) and the method detection limit (MDL) are given in the analytical method, which is described in reports on the investigation of the development of analytical methods for chemicals and adopted in the Initial Environmental Survey (hereafter, the Initial Environmental 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 2-2. The survey results are summarized as follows.

In surface water, 11 out of 22 target chemicals (groups) were detected.

- •[5] O-Ethyl O-4-nitrophenyl phenylphosphonothioate (synonym: EPN): 3 of 8 valid sites
- [7] 2,6-Xylenol: 2 of 6 valid sites
- [12] 3-(3,4-Dichlorophenyl)-1,1-dimethylurea (synonym: Diuron or DCMU): 9 of 10 valid sites
- [14] 2,4-Di-tert-butyl-6-(5-chloro-2*H*-1,2,3-benzotriazol-2-yl)phenol: 4 of 6 valid sites
- [19] O,O-Diethyl O-2-isopropyl-6-methyl-4-pyrimidinyl phosphorothioate (synonym: Diazinon): 7 of 10 valid sites
- [20] O,O-Dimethyl O-3-methyl-4-nitrophenyl phosphorothioate (synonym: Fenitrothion or MEP): 6 of 6 valid sites
- •[31] 2-(2*H*-1,2,3-Benzotriazol-2-yl)-4,6-di-*tert*-butylphenol: 2 of 6 valid sites
- [32] Methyl methacrylate: 1 of 7 valid sites
- [35] 2-sec-Butylphenyl N-methylcarbamate; (synonym: Fenobucarb or BPMC): 10 of 10 valid sites
- [37] Dimethyl 2,2-dichlorovinyl phosphate (synonym: Dichlorvos or DDVP): 6 of 8 valid sites
- [38] Tributyl phosphate: 10 of 19 valid sites

In sediment, 3 out of 7 target chemicals (groups) were detected.

- [14] 2,4-Di-*tert*-butyl-6-(5-chloro-2*H*-1,2,3-benzotriazol-2-yl)phenol: 6 of 6 valid sites
- [17] *N,N*-Dimethylformamide: 4 of 8 valid sites
- •[31] 2-(2*H*-1,2,3-Benzotriazol-2-yl)-4,6-di-*tert*-butylphenol: 6 of 6 valid sites

In wildlife (bivalves or fish), 5 out of the 11 target chemicals (groups) were detected.

- [12] 3-(3,4-Dichlorophenyl)-1,1-dimethylurea (synonym: Diuron or DCMU): 10 of valid sites
- [14] 2,4-Di-tert-butyl-6-(5-chloro-2H-1,2,3-benzotriazol-2-yl)phenol: 10 of 10 valid sites
- [18] Hydrogenated terphenyl: 2 of 10 valid sites
- •[27] Hydrazine: 9 of 10 valid sites
- •[31] 2-(2*H*-1,2,3-Benzotriazol-2-yl)-4,6-di-*tert*-butylphenol: 10 of 10 valid sites

In air, 6 out of the 7 target chemical were detected.

- [9] Isobutyl acetate: 4 of 7 valid sites
- •[13] 2,6-Dichlorobenzonitrile (synonym: Dichlobenil): 7 of 7 valid sites
- [22] Tetrahydrofuran: 3 of 7 valid sites
- [28] 1-Butanol: 5 of 7 valid sites
- [29] Furfural: 5 of 7 valid sites
- [33] 2-(1-Methylethoxy)ethanol: 3 of 7 valid sites

In food, 2 out of the 2 target chemical was detected.

- [27] Hydrazine: 146 samples of 178 valid samples
- [34] 2,3-Dihydro-2,2-dimethyl-7-benzo[b]furanyl *N*-methylcarbamate (synonym: Carbofuran): 14 samples of 178 valid samples

Table 2-2 Summary of the detection ranges and the detection limits in the Detailed Environmental Survey in FY 2006

Target chemicals		Surface water (ng/L)		Sediment (ng/g-dry)		Wildlife (ng/g-wet)		Air (ng/m³)		Food (ng/g-wet) or (ng/L)	
	27	Detection		Detection		Detection		Detection		Detection	
No.	Name	range and frequency	limit	range and frequency	limit	range and frequency	limit	range and frequency	limit	range and frequency	limit
[1]	2,2'-Azobisisobutyronitrile	nd 0/6	40							1	
[2]	2-Amino-4-[hydroxy(meth yl)phosphinoyl]butyric acid (synonym: Glufosinate)	nd 0/10	670								
[3]	m-Aminophenol	nd 0/7	7								
[4]	Aldicarb	nd 0/10	3	nd 0/6	0.17	nd 0/10	0.0016				
[5]	O-Ethyl O-4-nitrophenyl phenylphosphonothioate (synonym: EPN)	nd ~ 0.18 3/8	0.09								
[6-1]	N,N'-Ethylenebis(dithiocarba mamic acid) and its salt  Manganese N,N'-ethylenebis(dithiocarba mate) (synonym: Maneb)  Complex compounds of										
[6-2]	manganese  N,N'-ethylenebis(dithiocarba mate) and zinc  N,N'-ethylenebis(dithiocarba mate) (synonym: Mancozeb)  N,N'-Ethylenebis(thiocarbam	nd 0/7	30			nd 0/10	0.13				
[6-3]	oylthiozinc) bis( <i>N</i> , <i>N</i> -dimethyldithiocarba mate) (synonym: Polycarbamate)										
[7]	2,6-Xylenol	nd ~ 3.4 2/6	0.5								
[8]	Chlorobenzene			nd 0/6	0.3						
[9]	Isobutyl acetate							nd ~ 570 4/7	95		
[10]	Diisopropylnaphthalene	nd 0/4	0.4								
[11]	S-4-Chlorobenzyl N,N-diethylthiocarbamate (synonym: Thiobencarb)	nd 0/13	6								
[12]	3-(3,4-Dichlorophenyl)-1,1- dimethylurea (synonym: Diuron or DCMU)	nd ~ 230 9/10	0.6			nd ~ 0.20 10/10	0.0019				
[13]	2,6-Dichlorobenzonitrile (synonym: Dichlobenil)							0.10 ~ 0.76 7/7	0.04		
[14]	2,4-Di- <i>tert</i> -butyl-6-(5-chlor o-2 <i>H</i> -1,2,3-benzotriazol-2-yl)phenol	nd ~ 0.23 4/6	0.07	0.18 ~ 41 6/6	0.10	0.053 ~ 3.0 10/10	0.004				
[15]	N,N-Dimethyldithiocarbama mic acid and its salt										
[15- 1]	Zinc bis(N,N-dimethyldithiocarb amate) (synonym: Ziram) N,N'-Ethylenebis(thiocarbam oylthiozinc)	nd 0/7	50			nd 0/10	0.3				
[15- 2]	bis( <i>N</i> , <i>N</i> -dimethyldithiocarba mate) (synonym: Polycarbamate)										
[16]	N,N-Dimethyldodecylamine N-oxide			nd 0/4	0.8						
[17]	N,N-Dimethylformamide			nd ~ 18 4/8	1.4						

Target chemicals		Surface v (ng/L	.)	Sediment (ng/g-dry)		Wildlife (ng/g-wet)		Air (ng/m³)		Food (ng/g-wet) or (ng/L)	
No.	Name	Detecti range and		Detection range and		Detection range and	tion	Detecti range and	[	Dete range and	ction
110.	Name	frequency	limit	frequency	limit	frequency	limit	frequency	limit	frequency	limit
[18]	Hydrogenated terphenyl					nd ~ 0.81 2/10	0.52				
[19]	O,O-Diethyl O-2-isopropyl-6-methyl-4- pyrimidinyl phosphorothioate (synonym: Diazinon)	nd ~ 19 7/10	1								
[20]	O,O-Dimethyl O-3-methyl-4-nitrophenyl phosphorothioate (synonym: Fenitrothion or MEP)	nd ~ 4.8 6/6	0.011								
[21]	Tetrachloroisophthalonitrile (synonym: Chlorothalonil or TPN)	nd 0/8	0.5								
[22]	Tetrahydrofuran							nd ~ 260 3/7	60		
[23]	Trichloroacetaldehyde	nd 0/7	10								
[24]	Trichloronitromethane (synonym: Chloropicrin)					nd 0/10	0.3				
[25]	Nitrofen (synonym: NIP)							nd 0/6	0.7		
[26]	1,1-Bis( <i>tert</i> -butyldioxy)-3,3 ,5-trimethylcyclohexane					nd 0/10	0.03				
[27]	Hydrazine					nd ~ 95 9/10	1.2			diet nd ~ 0.80 (drinking water 0.77 ~ 2.7) 146/178	diet 0.0066 ~ 0.0095 (drinking water 0.68)
[28]	1-Butanol							nd ~ 1,400 5/7	60		
[29]	Furfural							nd ~ 85 5/7	40		
[30]	2-(2-Benzothiazolyloxy)- <i>N</i> -methylacetanilide (synonym: Mefenacet)	nd 0/13	25			nd 0/10	0.3		L		
[31]	22-(2 <i>H</i> -1,2,3-Benzotriazol-2-yl)-4,6-di- <i>tert</i> -butylpheno l	nd ~ 0.10 2/6	0.04	0.009 ~ 5.8 6/6	0.010	0.009 ~ 3.7 10/10	0.003				
[32]	Methyl methacrylate	nd ~ 15 1/7	8								
[33]	2-(1-Methylethoxy)ethanol							nd ~ 30 3/7	20		
[34]	2,3-Dihydro-2,2-dimethyl-7- benzo[b]furanyl <i>N</i> -methylcarbamate (synonym: Carbofuran)									diet nd ~ 0.12 (drinking water nd) 14/178	diet 0.0048 ~ 0.015 (drinking water 0.12)
[35]	2-sec-Butylphenyl N-methylcarbamate; (synonym: Fenobucarb or BPMC)	0.2 ~ 5.1 10/10	0.2								
[36]	<i>α</i> -Methylstyrene			nd 0/5	0.7						
[37]	Dimethyl 2,2-dichlorovinyl phosphate (synonym: Dichlorvos or DDVP)	nd ~ 20 6/8	0.3								

Target chemicals		Surface water		Sediment		Wildlife		Air		Food	
		(ng/L)		(ng/g-dry)		(ng/g-wet)		$(ng/m^3)$		(ng/g-wet) or (ng/L)	
		Detection		Detection		Detection		Detection		Detection	
No.	Name	range and frequency	limit	range and frequency	limit	range and frequency	limit	range and frequency	limit	range and frequency	limit
[38]	Tributyl phosphate	nd ~ 84 10/19	10								

- (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 were 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. In cases where a chemical is detected in one or more sameples from a site or an area, the site or area can be defined as one "detected site" or one "detected area".
- (Note 2) <u>Det</u>ection 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.
- (Note 4) is the sum value of detection limits of each congener and isomer, and therefore a detection range that does not exceed this value can be shown instead of "nd".
- (Note 5) Food is shown for diet (ng/g-wet) and drinking water (ng/L).