

**Henderson, Katie**

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**From:** Reiber, Loretta  
**Sent:** Tuesday, June 01, 2010 8:47 AM  
**To:** Henderson, Katie  
**Subject:** FW: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 1 of 4)

-----Original Message-----

**From:** Tillman.Michael@epamail.epa.gov [mailto:Tillman.Michael@epamail.epa.gov]  
**Sent:** Tuesday, August 04, 2009 9:38 AM  
**To:** bruce.fielding@la.gov  
**Cc:** Reiber, Loretta; Bailey, John; Baskin.Kilty@epamail.epa.gov  
**Subject:** Fw: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 1 of 4)

Good morning Bruce,

It was a pleasure speaking with you this morning, it has been a long while. As I mentioned, one of the permit writers with the state of Arkansas (Loretta Reiber) is wanting someone in your shop to look over the water quality screens they have performed for GP Crossett / Ouachita River Basin. At your earliest convenience, If you would please take a look at the attached information, for compliance with LA water quality standards that would be great.

I will be passing this along in 4 parts. Any questions, please contact Loretta directly at the phone number / email below. Thanks in advance for your time and assistance.

Mike Tillman (6WQ-PO)  
Arkansas / Texas - State NPDES Coord.  
U.S. EPA Region 6  
1445 Ross Ave.  
Dallas, TX 75202  
phn. 214-665-7531, fax. 214-665-2191

----- Forwarded by Michael Tillman/R6/USEPA/US on 08/04/2009 09:33 AM -----

**FW: GP Crossett, NPDES Permit No. AR0001210**

Reiber, Loretta

to: Michael Tillman

07/29/2009 02:39 PM

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Wouldn't let me send all at once so doing it in batches.

-----Original Message-----

**From:** Reiber, Loretta  
**Sent:** Wednesday, July 29, 2009 2:38 PM  
**To:** 'Tillman.Michael@epamail.epa.gov'  
**Subject:** GP Crossett, NPDES Permit No. AR0001210

6/1/2010

Mike,

Attached are the Priority Pollutant Scan calculations and ADEQ's review. ADEQ is requesting a review of these documents in order to confirm compliance with LDEQ's procedures at SMS2.

If you have any questions, please feel free to contact me at (501) 682-0612 or by e-mail at [reiber@adeq.state.ar.us](mailto:reiber@adeq.state.ar.us).

Loretta Reiber, P.E.  
Engineer, NPDES Permits

A. **Toxics Pollutants**

(1) Post Third Round Policy and Strategy

Section 101 of the Clean Water Act (CWA) states that "...it is the national policy that the discharge of toxic pollutants in toxic amounts be prohibited..." To insure that the CWA's prohibitions on toxic discharges are met, EPA has issued a "Policy for the Development of Water Quality-Based Permit Limitations by Toxic Pollutants"(49 FR 9016-9019, 3/9/84). In support of the national policy, Region 6 adopted the "Policy for post Third Round NPDES Permitting" and the "Post Third Round NPDES Permit Implementation Strategy" on October 1, 1992. The Regional policy and strategy are designed to insure that no source will be allowed to discharge any wastewater which (1) results in instream aquatic toxicity; (2) causes a violation of an applicable narrative or numerical State water quality standard resulting in non-conformance with the provisions of 40 CFR Part 122.44(d); (3) results in the endangerment of a drinking water supply; or (4) results in aquatic bioaccumulation which threatens human health.

(2) Implementation

The State of Arkansas is currently implementing EPA's Post Third-Round Policy in conformance with the EPA Regional strategy. The 5-year discharge permits contain technology-based effluent limitations reflecting the best controls available. Where these technology-based permit limits do not protect water quality or the designated uses, or where there are no applicable technology-based limits, additional water quality-based effluent limitations and/or conditions are included in the discharge permits. State narrative and numerical water quality standards from Regulation No. 2 are used in conjunction with EPA criteria and other available toxicity information to determine the adequacy of technology-based permit limits and the need for additional water quality-based controls.

(3) Priority Pollutant Scan

In accordance with the regional policy ADEQ has reviewed and evaluated the effluent in evaluating the potential toxicity of each analyzed pollutant:

- a. The results were evaluated and compared to EPA's Minimum Quantification Levels (MQLs) to determine the potential presence of a respective toxic pollutant. Those pollutants which are greater than or equal to the MQLs are determined to be reasonably present in the effluent and an evaluation of their potential toxicity is necessary.

- b. Those pollutants with one datum shown as "non-detect" (ND), providing the level of detection is equal to or lower than MQL are determined to be not potentially present in the effluent and eliminated from further evaluation.
- c. Those pollutants with a detectable value even if below the MQL are determined to be reasonably present in the effluent and an evaluation of their potential toxicity is necessary.
- d. For those pollutants with multiple data values and all values are determined to be non-detect, therefore no further evaluation is necessary. However, where data set includes some detectable concentrations and some values as ND, one-half of the detection level is used for those values below the level of detection to calculate the geometric mean of the data set.

The concentration of each pollutant after mixing with the receiving stream was compared to the applicable water quality standards as established in the Arkansas Water Quality Standards, Reg. No. 2 and with the aquatic toxicity, human health, and drinking water criteria obtained from the "Quality Criteria for Water, 1986 (Gold Book)". The manner in which the Instream Waste Concentrations are calculated may be found on page 2 of each of the attachments.

I. Aquatic Toxicity, Bioaccumulation, and Drinking Water

**Arkansas Requirements**

The flows (for acute, chronic, and bioaccumulation), TSS, hardness, etc. are based upon ADEQ's CPP.

<b>Outfall 001</b>		
Flow	52.4 MGD = 80.96 cfs	Application
7Q10	0 cfs	U.S.G.S.
TSS	5.5 mg/l	CPP, Section 5.24.3
Hardness as CaCO <sub>3</sub>	31 mg/l	CPP, Section 5.24.1
pH	7.01 s.u.	OUA008B
<b>Stream Monitoring Station (SMS2)</b>		
Flow	52.4 MGD = 80.96 cfs	Application
7Q10	1200 cfs	EPA*
TSS	5.5 mg/l	CPP, Section 5.24.3
Hardness as CaCO <sub>3</sub>	28 mg/l	CPP, Section 5.24.1
pH	7.01 s.u.	OUA008B

\*Letter dated July 3, 2001.

## Louisiana Requirements

The requirements of Louisiana are not applicable at Outfall 001 because of the distance from the outfall to the state line (over 10 stream miles). Also, effluent which is discharged through Outfall 001 is monitored at SMS2 when Mossy Lake is not flooded. SMS2 is a monitoring point located approximately 2.5 miles upstream of the Arkansas/Louisiana state line.

The flows (for acute, chronic, and bioaccumulation) are based upon the requirements of Title 33, Part IX, Subpart I, Section 1115, Table 2a.

Stream Monitoring Station (SMS2)		
Flow	52.4 MGD = 80.96 cfs	Application
7Q10	1200 cfs	EPA*
TSS	6 mg/l	E-mail**
Hardness as CaCO <sub>3</sub>	38.4 mg/l	E-mail**
pH	7.01 s.u.	OUA008B

\*Letter dated July 3, 2001.

\*\* These values were received via e-mail from Jeremy "Todd" Franklin of LDEQ on 06/16/2009.

### (4) Water Quality Standards for Metals and Cyanide

Standards for Chromium (VI), Mercury, Selenium, and Cyanide are expressed as a function of the pollutant's water-effect ratio (WER), while standards for cadmium, chromium (III), copper, lead, nickel, silver, and zinc are expressed as a function of the pollutant's water-effect ratio, and as a function of hardness.

The Water-effect ratio (WER) is assigned a value of 1.0 unless scientifically defensible study clearly demonstrates that a value less than 1.0 is necessary or a value greater than 1.0 is sufficient to fully protect the designated uses of the receiving stream from the toxic effects of the pollutant.

The WER approach compares bioavailability and toxicity of a specific pollutant in receiving water and in laboratory test water. It involves running toxicity tests for at least two species, measuring LC50 for the pollutant using the local receiving water collected from the site where the criterion is being implemented, and laboratory toxicity testing water made comparable to the site water in terms of chemical hardness. The ratio between site water and lab water LC50 is used to adjust the national acute and chronic criteria to site specific values.

### (5) Conversion of Dissolved Metals Criteria for Aquatic Life to Total Recoverable Metal

Metals criteria established in Regulation No. 2 for aquatic life protection are based on dissolved metals concentrations and hardness values (See Page 6 of Attachment 1). However, Federal Regulations cited at 40 CFR 122.45(c) require that effluent limitations for metals in discharge permits be expressed as total recoverable (See

Attachments 1, 2, and 3). Therefore a dissolved to the total recoverable metal conversion must be implemented. This involves determining a linear partition coefficient for the metal of concern and using this coefficient to determine the fraction of metal dissolved, so that the dissolved metal ambient criteria may be translated to a total effluent limit. The formula for converting dissolved metals to total recoverable metals for streams and lakes are provided in Section 5.25 of the CPP and Region 6 Implementation Guidance for Arkansas Water Quality Standards promulgated at 40 CFR Part 131.36.

(6) Comparison of the submitted information with the water quality standards and criteria

(a) Outfall 001

The following pollutants were determined to be present in the effluent from **Outfall 001** as reported by the permittee.

Pollutant	Concentration Reported, µg/l	MQL Required by ADEQ's CPP
Total Recoverable Chromium	4.3	10*
Hexavalent Chromium, Dissolved	4.3	10*
Total Recoverable Copper	6.79	0.5
Total Recoverable Lead	2.37	0.5
Total Recoverable Mercury	0.00833	0.005
Total Recoverable Nickel	7.07	0.5
Total Recoverable Selenium	2.22	5*
Total Recoverable Silver	3.58	0.5
Total Recoverable Thallium	2.96	0.5
Total Recoverable Zinc	373	20
Total Recoverable Phenols	0.0445	5*
Alpha-BHC	0.0501	0.05
Gamma-BHC	0.0642	0.05
Delta-BHC	0.0688	0.05
Endosulfan sulfate	0.0662	0.1
Endrin aldehyde	0.269	0.1

\*Actual detection level achieved was lower than what was required.

As indicated in the above table, ADEQ has determined from the information submitted by the permittee that the water quality standards for Total Recoverable Copper, Total Recoverable Lead, Total Recoverable Mercury, Total Recoverable Zinc, Alpha-BHC, Gamma-BHC, Delta-BHC, Endosulfan sulfate, and Endrin aldehyde and the Gold Book criteria for Total Recoverable Thallium are exceeded. Permit action will be taken for the parameters for which the permittee demonstrated reasonable potential for exceedances of the water quality standards and/or Gold Book Criteria (See Attachment 1).

(b) SMS2

The following pollutants were determined to be present at **SMS2** as reported by the permittee.

**Arkansas Standards**

Pollutant	Concentration Reported, µg/l	MQL Required by ADEQ's CPP
Total Recoverable Cadmium	1.17	1
Total Recoverable Chromium	5.66	10*
Hexavalent Chromium, Dissolved	5.66	10*
Total Recoverable Copper	11.7	0.5
Total Recoverable Lead	3.13	0.5
Total Recoverable Mercury	0.009	0.005
Total Recoverable Nickel	7.87	0.5
Total Recoverable Selenium	11.3	5
Total Recoverable Zinc	251	20
Bis(2-ethylhexyl) phthalate	<10	10*
Gamma-BHC	0.0969	0.05
Delta-BHC	<0.05	0.05*
Dieldrin	0.113	0.02
Alpha-endosulfan	0.0211	0.01

\*Actual detection level achieved was lower than what was required.

## Louisiana Standards

Pollutant	Concentration Reported, µg/l	MQL Required by LDEQ**
Total Recoverable Cadmium	1.17	1
Total Recoverable Chromium	5.66	10*
Hexavalent Chromium, Dissolved	5.66	10*
Total Recoverable Copper	11.7	10
Total Recoverable Lead	3.13	5
Total Recoverable Mercury	0.009	0.2
Total Recoverable Nickel	7.87	40
Total Recoverable Zinc	251	20
Gamma-BHC	0.0969	0.05
Dieldrin	0.113	0.1
Alpha-endosulfan	0.0211	0.1

\*Actual detection level achieved was lower than what was required.

\*\*Based on *Permitting Guidance Document for Implementing Louisiana Surface Water Quality Standards, Water Quality Management Plan, Volume 3*. Dated April 16, 2008 (Version 6). Obtained from LDEQ's web site on June 12, 2009.

\*\*\*WQS not listed in L.A.C. 33:IX:1113.

As indicated in the above tables, ADEQ has determined from the information submitted by the permittee that some water quality standards are exceeded while others are not. The effluent demonstrated reasonable potential for exceedances of Arkansas' water quality standards for Total Recoverable Selenium, Total Recoverable Zinc, and Dieldrin. The effluent also demonstrated reasonable potential for exceedances of Louisiana's water quality standards for Total Recoverable Zinc and gamma-BHC. Permit action will be taken for the parameters for which the permittee demonstrated reasonable potential for exceedances of the water quality standards (See Attachments 2 and 3). No Gold Book criteria were exceeded at SMS2.

### (c) Aquatic Toxicity

#### (i) Pollutants with numerical water quality standards

##### a. Outfall 001

ADEQ has determined from the information submitted by the permittee that there is a reasonable potential for the discharge to cause an instream excursion above the acute and/or chronic numeric standards as specified in the Arkansas Water Quality Standards, Reg. No. 2 (See Attachment 1).



ADEQ has identified the following toxicants in the discharge in amounts which could potentially have a toxic impact on the receiving stream:

**OUTFALL 001**

Chronic Aquatic Toxicity Results				
Pollutant	C <sub>e</sub> , µg/l	C <sub>e</sub> X 2.13	IWC, µg/l	AWQS, µg/l
Total Recoverable Copper	6.79	14.46	14.46	10.93
Total Recoverable Lead	2.37	5.05	5.05	3.40
Total Recoverable Mercury	0.00833	0.02	0.02	0.012
Total Recoverable Zinc	373	794.49	794.49	119.50
Alpha-BHC	0.0501	0.11	0.11	0.08
Gamma-BHC	0.0642	0.14	0.14	0.08
Delta-BHC	0.0688	0.15	0.15	0.08
Endosulfan Sulfate	0.0662	0.14	0.14	0.056
Endrin Aldehyde	0.269	0.57	0.57	0.0023

Acute Aquatic Toxicity Results				
Pollutant	C <sub>e</sub> , µg/l	C <sub>e</sub> X 2.13	IWC, µg/l	AWQS, µg/l
Total Recoverable Silver	3.58	7.63	7.63	1.51
Total Recoverable Zinc	373	794.49	794.49	130.87
Endrin Aldehyde	0.269	0.57	0.57	0.18

b. SMS2

ADEQ has determined from the information submitted by the permittee that there is a reasonable potential for the discharge to cause an instream excursion above the acute and/or chronic numeric standards as specified in the Arkansas Water Quality Standards, Reg. No. 2 and/or in Louisiana's Water Quality Regulations at L.A.C. 33:IX:1113 (See Attachments 2 and 3).

ADEQ has identified the following toxicants in the discharge in amounts which could potentially have a toxic impact on the receiving stream:

Chronic Aquatic Toxicity Results						
Pollutant	C <sub>e</sub> , µg/l	C <sub>e</sub> X 2.13	AR IWC, µg/l	AR WQS, µg/l	LA IWC, µg/l	LA WQS, µg/l
Dieldrin	0.113	0.241	0.05	0.019	**	**
Total Recoverable Selenium	11.3	24.07	5.11	5	*	*
Total Recoverable Zinc	251	534.63	113.62	109.63	**	**
Gamma-BHC	0.0969	0.206	***	***	0.35	0.21

\*WQS not listed in L.A.C. 33:IX:1113.

\*\*Reasonable potential only demonstrated based upon Louisiana's requirements.

\*\*\*Reasonable potential only demonstrated based upon Arkansas' requirements.

Acute Aquatic Toxicity Results						
Pollutant	C <sub>e</sub> , µg/l	C <sub>e</sub> X 2.13	AR IWC, µg/l	AR WQS, µg/l	LA IWC, µg/l	LA WQS, µg/l
Total Recoverable Zinc	251	534.63	282.97	120.05	359.02	159.70

IWC's have been calculated in the manner described on page 2 of the attachments.

#### c. Permit Action

Under Federal Regulation 40 CFR Part 122.44(d), as adopted by Regulation No. 6, if a discharge poses the reasonable potential to cause or contribute to an exceedance above a water quality standard, the permit must contain an effluent limitation for that pollutant. Effluent limitations for the toxicants listed above have been derived in a manner consistent with the Technical Support Document (TSD) for Water Quality-based Toxics Control (EPA, March 1991), the State's implementations procedures, and 40 CFR Part 122.45(c).

#### Permit Limit Determination

The instream waste load allocation (WLA), which is the level of effluent concentration that would comply with the water quality standard (WQS) of the receiving stream, is calculated for both chronic and acute WLA using the following equations:

$$WLA_c = (WQS \times (Q_d + Q_b) - Q_b \times C_b) / Q_d$$

Where:

WLA<sub>c</sub> = chronic waste load allocation (µg/l)

$Q_d$  = discharge flow (cfs)  
 $Q_b$  = 0.67 X 7Q10 (cfs) @ **Outfall 001**  
 $Q_b$  = 0.25 X 7Q10 (cfs) @ **SMS2 for AR WQS**  
 $Q_b$  = 0.33 X 7Q10 (cfs) @ **SMS2 for LA WQS**  
 $C_b$  = background concentration ( $\mu\text{g/l}$ )  
 WQS = chronic aquatic toxicity standards ( $\mu\text{g/l}$ )

and;

$$WLA_a = (WQS \times (Q_d + Q_b) - Q_b \times C_b) / Q_d$$

Where:

$WLA_a$  = acute waste load allocation ( $\mu\text{g/l}$ )  
 $Q_d$  = discharge flow (cfs)  
 $Q_b$  = 0.33 X 7Q10 (cfs) @ **Outfall 001**  
 $Q_b$  = 0.06 X 7Q10 (cfs) @ **SMS2 for AR WQS**  
 $Q_b$  = 0.033 X 7Q10 (cfs) @ **SMS2 for LA WQS**  
 $C_b$  = background concentration ( $\mu\text{g/l}$ )  
 WQS = acute aquatic toxicity standards ( $\mu\text{g/l}$ )

The long term average (LTA) effluent concentration is then calculated based on the chronic and acute WLA as follows:

$$LTA_c = 0.72 \times WLA_c$$

$$LTA_a = 0.57 \times WLA_a$$

The lowest of these two (2) values is selected as being the limiting LTA. The limiting LTA is then used to calculate the monthly average (AML) and daily maximum (DML) for the final limits. AML and DML are calculated as follows:

$$AML = 1.55 \times \text{Limiting LTA}$$

$$DML = 3.11 \times \text{Limiting LTA}$$

Limits included in the permit are as follows:

### **OUTFALL 001**

See Attachment 1 regarding the calculations for the following permit limits.

Arkansas Numerical Aquatic Toxicity Limits		
Parameter	AML*, $\mu\text{g/l}$	DML*, $\mu\text{g/l}$
Total Recoverable Copper	12.20	24.48

Arkansas Numerical Aquatic Toxicity Limits		
Parameter	AML*, µg/l	DML*, µg/l
Total Recoverable Lead	3.80	7.62
Total Recoverable Mercury	0.012	0.012
Total Recoverable Zinc	115.62	231.99
Alpha-BHC	0.04	0.12
Gamma-BHC	0.09	0.18
Delta-BHC	0.09	0.18
Endosulfan Sulfate	0.06	0.13
Endrin Aldehyde	0.003	0.005

## SMS2

See Attachment 2 regarding the calculations for the limits on Total Recoverable Selenium, Total Recoverable Zinc, and Dieldrin. The permittee did not demonstrate reasonable potential for water quality violations due to Total Recoverable Selenium and Dieldrin in Louisiana. Therefore, those permit limits were based on ADEQ's permitting procedures.

The reach of the Ouachita River which receives the effluent from this facility (Reach #002 in H.U.C. 08040202) is on the 303(d) due to Zinc. The permittee demonstrated reasonable potential for water quality violations due to the levels of Total Recoverable Zinc. Permit limits were determined using both ADEQ's and LDEQ's permitting procedures. The more stringent limits were those calculated using ADEQ's permitting procedures. Therefore, those limits were placed in the permit.

The permittee did not demonstrate reasonable potential for water quality violations due to Total Recoverable Copper in either Arkansas or Louisiana. However, the reach of the Ouachita River which receives the effluent from this facility is on the 303(d) list due to Copper. Therefore, permit limits based on ADEQ's permitting procedures have been included in the permit.

The Department will reopen the permit to include any TMDLs which are finalized during the term of the permit.

See Attachment 3 regarding the calculations for the limits on Gamma-BHC. The permittee did not show reasonable potential for water quality

violations due to Gamma-BHC in Arkansas. Therefore, the Gamma-BHC limits were based upon LDEQ's permitting procedures.

Numerical Aquatic Toxicity Limits		
Parameter	AML*, µg/l	DML*, µg/l
Gamma-BHC	1.381	2.770
Total Recoverable Copper	22.43	45.00
Total Recoverable Selenium	26.26	52.68
Total Recoverable Zinc	200.40	402.09
Dieldrin	0.01	0.02

(ii) Pollutants without applicable water quality standards

ADEQ has determined from the information submitted by the permittee that there is not a reasonable potential for the discharge to cause an instream excursion above the acute and/or chronic criteria as specified in the Gold Book (See Attachments 1, 2, and 3).

b. Human Health (Bioaccumulation) Limits

i. Pollutants with numerical water quality standards

ADEQ has determined from the information submitted by the permittee that there is not a reasonable potential for the discharge to cause an instream excursion above the state numeric bioaccumulation standards as specified in Reg. 2.508 and LDEQ's water quality regulations.

ii. Pollutants without applicable water quality standards

ADEQ has determined from the information submitted by the permittee that there is reasonable potential for the discharge to cause exceedence of bioaccumulation criterion as specified in the Gold Book (Quality Criteria for Water 1986) for only **Total Recoverable Thallium at Outfall 001**. The results of the analysis are as follows:

Bioaccumulation Criterion Results				
Pollutant	C <sub>e</sub> , µg/l	C <sub>e</sub> X 2.13	IWC, µg/l	GB, µg/l
Total Recoverable Thallium	2.96	6.3048	6.3048	6.3

IWC's have been calculated in the manner described on page 2 of the attachments.

Since the Arkansas Water Quality Standards have not been established for those parameters listed above, no permit limitations have been placed in the draft permit. However, monitoring and reporting is required to confirm that the pollutant is present at the levels reported by the permittee. The permit may be reopened to require effluent limitations, additional testing, and/or other appropriate actions.

iii. Drinking Water Supply Protection

ADEQ has determined from the information submitted by the permittee that there is not a reasonable potential for the discharge to cause an instream excursion above the drinking water criteria as specified in the Gold Book.

**Henderson, Katie**

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**From:** Reiber, Loretta  
**Sent:** Tuesday, June 01, 2010 8:47 AM  
**To:** Henderson, Katie  
**Subject:** FW: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 2 of 4)

-----Original Message-----

**From:** Tillman.Michael@epamail.epa.gov [mailto:Tillman.Michael@epamail.epa.gov]  
**Sent:** Tuesday, August 04, 2009 9:40 AM  
**To:** bruce.fielding@la.gov  
**Cc:** Reiber, Loretta; Bailey, John; Baskin.Kilty@epamail.epa.gov  
**Subject:** Fw: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 2 of 4)

----- Forwarded by Michael Tillman/R6/USEPA/US on 08/04/2009 09:48 AM -----

**FW: GP Crossett, NPDES Permit No. AR0001210**



Reiber, Loretta

to: Michael Tillman

07/29/2009 02:41 PM

-----Original Message-----

**From:** Reiber, Loretta  
**Sent:** Wednesday, July 29, 2009 2:38 PM  
**To:** 'Tillman.Michael@epamail.epa.gov'  
**Subject:** GP Crossett, NPDES Permit No. AR0001210

Mike,

Attached are the Priority Pollutant Scan calculations and ADEQ's review. ADEQ is requesting a review of these documents in order to confirm compliance with LDEQ's procedures at SMS2.

If you have any questions, please feel free to contact me at (501) 682-0612 or by e-mail at [reiber@adeq.state.ar.us](mailto:reiber@adeq.state.ar.us).

Loretta Reiber, P.E.  
Engineer, NPDES Permits

A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
1															
2															
3															
4	STEP 1:	INPUT TWO LETTER CODE FOR ECOREGION (Use Code at Right)													
5		Basin Name													
6															
7	FACILITY														
8															
9	Permittee														
10	NPDES Permit No.														
11	Outfall No(s)														
12	Plant Effluent Flow (MGD)														
13	Plant Effluent Flow (cfs)														
14															
15	RECEIVING STREAM														
16															
17	Is this a large river? (see list at right)/enter '1' if yes, '0' if no, make entry as a number														
18	Name of Receiving Stream														
19	Waterbody Segment Code No.														
20	Is this a lake or reservoir? (enter '1' if yes, '0' = no; make entry as a number)														
21	Second Enter 7Q10 in Cell H31														
22	(Reserved)														
23	(Reserved)														
24	(Reserved)														
25	(Reserved)														
26	(Reserved)														
27	(Reserved)														
28															
29	Ecotone TSS (mg/l) (For Large River, See List to Right)														
30	Ecotone Hardness (mg/l)														
31	Enter 7Q10 (cfs) as the Critical Flow														
32	Long Term Ave / Harmonic Mean Flow (cfs)														
33	Using Diffusers (Yes/No)														
34	pH (Avg)														
35	Percent (%) of Critical Flow for Chronic Criteria														
36	Percent (%) of Critical Flow for Acute Criteria														
37	Water Effect Ratio (WER)														
38	Ave Monthly Limit LTA Multiplier (Ref. page 103 TSD for WQ-Based Toxics Control)														
39	Max Daily Limit LTA Multiplier (Ref. page 103 TSD for WQ-Based Toxics Control)														

Ouachita Mts. Eco (OM) = 2.0 mg/l  
 Ozark Highlands Eco (OH) = 2.5 mg/l  
 Boston Mts. Eco (BM) = 1.3 mg/l  
 Ark River Valley Eco (AV) = 3.0 mg/l  
 Gulf Coastal Eco (GC) = 5.5 mg/l  
 Delta Ecotone (DL) = 8.0 mg/l  
 Ouachita (Above Caddo River) = 2.0 mg/l  
 Ouachita (Below Caddo River) = 5.5 mg/l  
 Red River = 33.0 mg/l  
 Arkansas (Ft. Smith to Dardanelle Dam) = 12.0 mg/l  
 Arkansas (Dardanelle Dam to Terry L&L) = 10.5 mg/l  
 Arkansas (Terry L&L to L&D No. 5) = 8.3 mg/l  
 Arkansas (L&D No. 5 to Mouth) = 9.0 mg/l  
 White (Above Beaver Lake) = 2.5 mg/l  
 White (Below Bull Shoals to Black Riv) = 3.3 mg/l  
 White (From Black River to Mouth) = 18.5 mg/l  
 St. Francis River = 18.0 mg/l  
 Ouachita (Above Caddo River) = 2.0 mg/l  
 Ouachita (Below Caddo River) = 5.5 mg/l  
 Red River = 33.0 mg/l  
 Total Hardness for:  
 Arkansas River = 125 mg/l  
 Ouachita River = 28 mg/l  
 White River = 116 mg/l  
 Gulf Coastal = 31 mg/l  
 Ozark Highlands = 148 mg/l  
 Boston Mount = 25 mg/l  
 Large Rivers  
 Mississippi River, Arkansas River, Red River  
 White (Below confluence with Black River)  
 Ouachita (Below confluence with Little Miss. River)  
 For industrial and federal facility, use the highest monthly average flow for the past 24 months. For POTWs, use the design flow.  
 #VALUE! => No violation or Not Applicable  
 9999999.00 => No EPA/ADEQ Guideline



	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
40																
41	STEP 2:	INPUT AMBIENT AND EFFLUENT DATA														
42		CALCULATE IN-STREAM WASTE CONCENTRATIONS														
43		For less than 20 data points enter geometric mean concentration as micro-gram per liter (ug/l or ppb)														
44	DATA INPUT	For 20 or more data points in set enter highest concentration as micro-gram per liter (ug/l or ppb)														
45																
46		Effluent value reported as "< detection level" (DL) but the DL is greater than MDL, the 1/2 DL is used.														
47		Effluent value reported as "< detection level" (DL) and the DL is smaller than MDL, "0" is used.														
48		If a firm value is reported, even less than MDL, the reported value is used														
49																
50																
51		The following formulae is used to calculate the instream Waste Concentration (IWC)														
52		(Please refer to CPP for detail)														
53		$IWC = [(F \cdot Q_a \cdot C_b) + (Q_e \cdot 2.13 \cdot C_e)] / (F \cdot Q_a + Q_e)$														
54		Where:														
55		IWC = Instream Waste Concentration														
56		F = Fraction of stream allowed for mixing														
57		C <sub>e</sub> = Reported concentration in effluent														
58		C <sub>b</sub> = Ambient stream concentration upstream of discharge														
59		Q <sub>e</sub> = Plant effluent flow														
60		Q <sub>b</sub> = Critical low flow of stream at discharge point expressed as the 7Q10 or harmonic mean flow for human health criteria														
61		Upstream Flow (Q <sub>b</sub> ) = (% of 7Q10) X 7Q10 for Chronic and Acute														
62																
63		The following formulae convert metals reported in total form to dissolved form if criteria are in dissolved form														
64																
65		$K_p = K_{pp} \cdot (TSS^{**a})$														
66		$C/C_i = 1 / (1 + K_p \cdot TSS \cdot 10^{-6})$														
67		Total Metal Criteria (Ct) = Cr / (C/Ci)														
68																
69		*Stream Linear Partition Coefficient (Insert "Dissolved" Conc in Column B to convert to Lake Linear Partition Coefficient														
70	Total Metals	K <sub>pp</sub>	alpha (a)	K <sub>p</sub>	C/Ci	Total Value	K <sub>pp</sub>	alpha (a)	K <sub>p</sub>	C/Ci	Total Value	K <sub>pp</sub>	alpha (a)	K <sub>p</sub>	C/Ci	Total Value
71	Arsenic	4800000	-0.73	129774.936	0.562224279	0.00	4800000.00	-0.73	129774.94	0.5622243	0	4800000.00	-0.73	129774.94	0.5622243	0
72	Cadmium	40000000	-1.13	528140.021	0.239874874	0.00	35200000.00	-0.92	677085.50	0.1975304	0	21700000.00	-0.27	1337700.52	0.1107886	0
73	Chromium(3)	33600000	-0.93	634831.714	0.207943859	0.00	28500000.00	-0.9	568209.82	0.2267955	0	20400000.00	-0.53	789241.87	0.1743542	0
74	Copper	10400000	-0.74	276185.843	0.376348023	0.00	19700000.00	-1.17	242119.22	0.4077114	0	22100000.00	-0.76	566235.80	0.2274063	0
75	Lead	28000000	-0.8	687785.571	0.199731923	0.00	33400000.00	-0.57	967651.25	0.1443854	0	24000000.00	-1.03	379066.54	0.3053995	0
76	Mercury	29000000	-1.14	376101.939	0.307067625	0.00	12500000	-0.7	356618.721	0.318500517	0.00	24000000	-1.03	379066.542	0.305399532	0.00
77	Nickel	49000000	-0.57	176461.46	0.485727208	0.00	24000000	-1.03	379066.542	0.305399532	0.00	24000000	-1.03	379066.54	0.3053995	0
78	Zinc	12500000	-0.7	356618.721	0.318500517	0.00	24000000	-1.03	379066.542	0.305399532	0.00	24000000	-1.03	379066.54	0.3053995	0
79	Silver	24000000	-1.03	379066.542	0.305399532	0.00	24000000	-1.03	379066.542	0.305399532	0.00	24000000	-1.03	379066.54	0.3053995	0
80		*Note: Use this section to compare lab concentrations shown as "Dissolved" vs "Total"														
81																



A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
132 POLLUTANTS		Number of Data points	MQL	EPA Statistical Factor	Background Conc. Cb ug/l	Effluent Conc. Ce ug/l	Domestic Supply IWC ug/l	Acute Aquatic IWC ug/l	Chronic Aquatic IWC ug/l	Bioacc. IWC ug/l	Domestic Criteria ug/l	Arkansas Acute Criteria ug/l	Arkansas Chronic Aquatic ug/l	Arkansas Bioacc. ug/l	
133															
<b>137 METALS AND CYANIDE</b>															
138	1. Antimony Total	1	60	2.13	0	0	0.00	0.00	0.00	0.00	#####	9000.00	1600.00	10	
139	2. Arsenic Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	50	640.31	337.94	1.4	
140	3. Beryllium Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	#####	130.00	5.30	4	
141	4. Cadmium Total	1	1	2.13	0	1.17	0.16	1.67	0.42	0.16	10	5.45	2.12	10.00	
142	5. Chromium (Tr)	1	10	2.13	0	566	0.76	8.10	2.05	0.76	50	1205.00	390.89	50.00	
143	7. Chromium (hex)	1	10	2.13	0	566	0.76	8.10	2.05	0.76	50	1205.00	390.89	50.00	
144	8. Copper Total	1	0.5	2.13	0	11.7	1.58	16.74	4.23	1.58	#####	19.89	14.42	1000.00	
145	9. Lead Total	1	0.5	2.13	0	3.13	0.42	4.48	1.13	0.42	50	112.47	4.38	50.00	
146	10. Mercury Total	1	0.005	2.13	0	0.009	0.00	0.01	0.00	0.00	2	6.64	0.012	2	
147	12. Nickel Total	1	0.5	2.13	0	7.87	1.06	11.26	2.85	1.06	#####	1296.68	144.01	4600	
148	13. Selenium Total	1	5	2.13	0	11.3	1.52	16.16	4.09	1.52	10	20.00	5.00	#####	
149	14. Silver Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	50	2.18	#####	6.3	
150	15. Thallium Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	#####	1400.00	#####	#####	
151	16. Zinc Total	1	20	2.13	0	251	33.79	359.02	90.75	33.79	#####	159.70	145.83	#####	
152	129. Phenols Total	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	700.00	350.00	#####	
153	17. Cyanide Total	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	45.90	5.4	663.8	
<b>155 DIOXIN</b>															
156	18. 2-3-7-8-TCDD	1	0.00001	2.13	0	0	0.00	0.00	0.00	0.00	#####	0.01	1.00E+07	7.10E-07	
<b>159 VOLATILE COMPOUNDS</b>															
160	19. Acrolein	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	68	21	780	
161	20. Acrylonitrile	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	7550	2600	6.6	
162	21. Benzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	2249	1125	0.58	
163	22. Bromoform	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	2930	1465.00	3.9	
164	23. Carbon Tetrach	1	2	2.13	0	0	0.00	0.00	0.00	0.00	5	2730	1365.00	0.22	
165	24. Chlorobenzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	250	50	2.10E+04	
166	25. Chlorobromomethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	340	
167	26. Chloroethane	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	1.00E+07	
168	27. 2-Chloroethylvinyl ether	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	1.00E+07	
169	28. Chloroform	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	2890	1445	5.3	
170	29. Dichlorobromomethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	220	
171	30. 1-1-Dichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	7	9999999.00	#####	#####	
172	31. 1-2-Dichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	11800	5900	0.36	
173	32. 1-1-Dichloroethylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1160	580.00	0.05	
174	33. 1,2-Dichloropropane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	23000	5700	#####	
175	34. 1,5-Dienorpropylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	6060	244	1700	
176	35. Ethylbenzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	3200	1600.00	29000	
177	37. Methyl Chloride	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	55000.00	275000.00	#####	
178	36. Methyl bromide	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	4000	
179	38. Methylene chloride	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	19300.00	9650.00	4.4	
180	39. 1-1-2,2-Tetrachloroetha	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	932	466	0.16	
181	40. Tetrachloroethylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1290	645	0.65	
182	41. Toluene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1270	635.00	6.10E+03	
183	42. 1,2-trans-dichloroethyl	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
184	44. 1-1-2-Trichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1800	900	0.56	
185	43. 1-1-1-Trichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	200	5280	2640.00	200.00	
186	45. Trichloroethylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	3900	1950	2.8	
187	46. Vinyl Chloride	1	10	2.13	0	0	0.00	0.00	0.00	0.00	2	9999999.00	#####	0.0237	

NO	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
						Ambient Background Conc. Cb	Effluent Conc. Ce	Domestic I/WC Supply	Acute I/WC	Chronic Aqueatic I/WC	Human Health I/WC	Domestic Criteria ug/l	Acute Aqueatic Criteria ug/l	Chronic Aqueatic Criteria ug/l	Human Health Criteria ug/l	
189	<b>ACID COMPOUNDS</b>															
190	47	2-Chlorophenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	258	129.00	0.10	
191	48	2,4-Dichlorophenol	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	202	101	0.30	
192	49	2,4-Dimethylphenol	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
193	50	4,6-Dinitro-o-Cresol	1	50	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	765	
194	51	2,4-Dinitrophenol	1	50	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14000	
195	52	53-Nitrophenols	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	230	150	#####	
196	54	4-Chloro-3-methylphenol	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	30	#####	#####	
197	55	Pentachlorophenol	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9.16	5.78	82	
198	56	Phenol	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	10200	2560	4600000	
199	57	2,4,6-Trichlorophenol	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	65	
201	<b>BASE/NEUTRAL COMPOUNDS</b>															
202	58	Acenaphthene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	1700	520	#####	
203	59	Acenaphthylene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
204	60	Anthracene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	110000	
205	61	Benzidine	1	50	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	250	125.00	8.00E-05	
207	62	Benzol(a) anthracene	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	
208	63	Benzol(a) pyrene	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	
209	64	3,4-benzofluoranthene	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
210	65	Benzo(g,h,i)perylene	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
211	66	Benzo(k) fluoranthene	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	
212	67	Bis(2-chloroethoxy)meth	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14	
213	68	Bis(2-chloroethyl) Ether	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
214	69	Bis(2-Chloroisopropyl) e	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	1.70E+05	
215	70	Bis(2-ethylhexyl)phthalate	1	10	2.13	0	2.87	0.39	4.11	1.04	0.39	#####	9999999.00	#####	59	
216	71	4-Bromophenyl phenyle	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
217	72	Butylbenzyl phthalate	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	1600	#####	#####	
218	73	2-chloronaphthalene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	
219	74	4-chlorophenyl phenyle	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	
220	75	Chrysene	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	1120	763	2600	
221	76	Dibenz(a,h)anthracene	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.77	
222	77-79	Dichlorobenzene(1,2	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	1.20E+05	
223	80	3,3'-Dichlorobenzidine	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	2.90E+06	
224	81	Diethyl Phthalate	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	1.20E+04	
225	82	Dimethyl phthalate	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	270	#####	5.4	
226	83	D-n-Butyl phthalate	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	3980	#####	370	
227	84	2,4-Dinitrotoluene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
228	85	2,6-Dinitrotoluene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	330	230	91	
229	86	D-n-octyl phthalate	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
230	87	1,2-diphenylhydrazine	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	270	#####	#####	
231	88	Fluoranthene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14000	
232	89	Fluorene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.00025	
233	90	Hexachlorobenzene	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	5.10	1.02	0.09	
234	91	Hexachlorobutadiene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	7	5.2	1.74E+04	
235	92	Hexachlorocyclopentadi	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	980	540	89	
236	93	Hexachloroethane	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	2	0.08	#####	
237	94	Hexachlorocyclohexane	1	5	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	
238	94	Indeno(1,2,3-cd)pyrene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	117000	#####	6000	
239	95	Isocourene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	2300	620	#####	
240	96	Naphthalene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	27000	#####	1900	
241	97	Nitrobenzene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	81	
242	98	N-nitrosodimethylamine	1	50	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
243	99	N-nitrosod-n-propylamir	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	160	
244	100	N-nitrosodiphenylamir	1	20	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
245	101	Phenanthrene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
246	103	1,2,4-trichlorobenzene	1	10	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
247																
248	<b>PESTICIDES</b>															
249	104	Aldrin	1	0.01	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	3.00	#####	0.00004	
250	105	Alpha-BHC	1	0.05	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	2.00	0.08	0.0373	
251	106	Beta-BHC	1	0.05	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	2.00	0.08	0.46	
252	107	Gamma-BHC	1	0.05	2.13	0	0.969	0.13	1.39	0.35	0.13	#####	5.30	0.21	0.11	
253	108	Delta-BHC	1	0.05	2.13	0	0.0407	0.01	0.06	0.01	0.01	#####	2.00	0.08	#####	
254	109	Chlordane	1	0.2	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	2.40	0.0043	0.00019	
255	110	4,4'-DDE	1	0.02	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	52.50	10.5	0.00019	
256	111	4,4'-DDE	1	0.1	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	0.03	0.006	0.00027	
257	112	4,4'-DDD	1	0.1	2.13	0	0.00	0.00	0.00	0.00	0.00	#####	0.03	0.006	0.00027	





**Henderson, Katie**

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**From:** Reiber, Loretta  
**Sent:** Tuesday, June 01, 2010 8:48 AM  
**To:** Henderson, Katie  
**Subject:** FW: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 3 of 4)

-----Original Message-----

**From:** Tillman.Michael@epamail.epa.gov [mailto:Tillman.Michael@epamail.epa.gov]  
**Sent:** Tuesday, August 04, 2009 9:42 AM  
**To:** bruce.fielding@la.gov  
**Cc:** Reiber, Loretta; Bailey, John; Baskin.Kilty@epamail.epa.gov  
**Subject:** Fw: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 3 of 4)

---- Forwarded by Michael Tillman/R6/USEPA/US on 08/04/2009 09:51 AM ----

**FW: GP Crossett, NPDES Permit No. AR0001210**



Reiber, Loretta

to: Michael Tillman

07/29/2009 02:41 PM

-----Original Message-----

**From:** Reiber, Loretta  
**Sent:** Wednesday, July 29, 2009 2:38 PM  
**To:** 'Tillman.Michael@epamail.epa.gov'  
**Subject:** GP Crossett, NPDES Permit No. AR0001210

Mike,

Attached are the Priority Pollutant Scan calculations and ADEQ's review. ADEQ is requesting a review of these documents in order to confirm compliance with LDEQ's procedures at SMS2.

If you have any questions, please feel free to contact me at (501) 682-0612 or by e-mail at [reiber@adeq.state.ar.us](mailto:reiber@adeq.state.ar.us).

Loretta Reiber, P.E.  
Engineer, NPDES Permits





	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
40																
41	STEP 2:															
42	INPUT AMBIENT AND EFFLUENT DATA															
43	CALCULATE IN-STREAM WASTE CONCENTRATIONS															
44	DATA INPUT															
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\*Please Use this section to convert lab concentrations shown as mg/L to ug/L



	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
132	POLLUTANTS		Number of Data points	MQL ug/l	EPA Statistical Factor	Background Conc. ug/l	Effluent Conc. ug/l	Domestic Supply IWC ug/l	Acute Aquatic IWC ug/l	Chronic Aquatic IWC ug/l	Bioacc. IWC ug/l	Domestic Criteria ug/l	Arkansas Acute Aquatic Criteria ug/l	Arkansas Chronic Aquatic ug/l	Arkansas Bioacc. ug/l	
133																
137	METALS AND CYANIDE															
138	1. Antimony Total		1	60	2.13	0	0	0.00	0.00	0.00	0.00	#####	9000.00	1600.00	4300	
139	2. Arsenic Total		1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	50	633.81	334.51	1.4	
140	3. Beryllium Total		1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	#####	130.00	5.30	4	
141	4. Cadmium Total		1	1	2.13	0	0	0.00	0.00	0.00	0.00	10	4.37	1.82	#####	
142	6. Chromium (Tr)		1	10	2.13	0	4.3	9.16	9.16	9.16	9.16	50	1006.35	326.45	#####	
143	7. Chromium (Hex)		1	10	2.13	0	4.3	9.16	9.16	9.16	9.16	50	15.71	10.58	#####	
144	8. Copper Total		1	0.5	2.13	0	6.79	14.46	14.46	14.46	14.46	#####	14.79	10.93	#####	
145	9. Lead Total		1	0.5	2.13	0	2.37	5.05	5.05	5.05	5.05	50	87.29	3.40	#####	
146	10. Mercury Total		1	0.005	2.13	0	0.00833	0.02	0.02	0.02	0.02	2	6.70	0.012	0.15	
147	12. Nickel Total		1	0.5	2.13	0	7.07	15.06	15.06	15.06	15.06	#####	101.45	117.88	4600	
148	13. Selenium Total		1	5	2.13	0	2.22	4.73	4.73	4.73	4.73	10	20.00	5.00	#####	
149	14. Silver Total		1	0.5	2.13	0	3.58	7.63	7.63	7.63	7.63	50	1.51	#####	6.3	
150	15. Thallium Total		1	0.5	2.13	0	2.96	6.3048	6.3048	6.3048	6.3048	#####	1400.00	#####	#####	
151	16. Zinc Total		1	20	2.13	0	373	794.49	794.49	794.49	794.49	#####	130.87	119.50	#####	
152	129. Phenols, Total		1	5	2.13	0	0.0445	0.09	0.09	0.09	0.09	#####	9999999.00	#####	#####	
153	17. Cyanide Total		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	22.36	5.2	220000	
156	DIOXIN															
157	18. 2-3,7-8-TCDD		1	0.00001	2.13	0	0	0.00	0.00	0.00	0.00	#####	0.01	1.00E+07	1.00E-06	
159	VOLATILE COMPOUNDS															
160	19. Acrolein		1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	68	21	780	
161	20. Acrylonitrile		1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	7550	2600	6.6	
162	21. Benzene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	5300	9999999	710	
163	22. Bromoform		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	3600	
164	23. Carbon Tetrach		1	2	2.13	0	0	0.00	0.00	0.00	0.00	5	35200	#####	44	
165	24. Chlorobenzene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	250	50	2.10E+04	
166	25. Chlorobromomethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	340	
167	26. Chloroethane		1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	1.00E+07	
168	27. 2-Chloroethylvinyl ether		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	1.00E+07	
169	28. Chloroform		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	28900	1240	4700	
170	29. Dichlorobromomethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	7	9999999.00	#####	220	
171	30. 1,1-Dichloroethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
172	31. 1,2-Dichloroethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	11800	20000	990	
173	32. 1,1-Dichloroethylene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	11600	#####	32	
174	33. 1,2-Dichloropropane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	23000	5700	#####	
175	34. 1,3-Dichloropropylene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	6060	244	1700	
176	35. Ethylbenzene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	32000	#####	29000	
177	37. Methyl Chloride		1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
178	38. Methyl bromide		1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	4000	
179	38. Methylene chloride		1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	16000	
180	39. 1-1,2,2-Tetrachloroethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9320	2400	110	
181	40. Tetrachloroethylene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	5280	840	88.5	
182	41. Toluene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	17500	#####	2.00E+05	
183	42. 1,2-trans-dichloroethylene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
184	44. 1-1,2-Trichloroethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	18000	9400	420	
185	43. 1-1,1-Trichloroethane		1	10	2.13	0	0	0.00	0.00	0.00	0.00	200	18000	#####	#####	
186	45. Trichloroethylene		1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	45000	21900	810	
187	46. Vinyl Chloride		1	10	2.13	0	0	0.00	0.00	0.00	0.00	2	9999999.00	#####	5250	

	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
						Ambient Background Conc.	Effluent Conc.	Domestic Supply	Acute IWC	Chronic IWC	Human Health IWC	Domestic Criteria ug/l	Acute Aquatic Criteria ug/l	Chronic Aquatic Criteria ug/l	Human Health Criteria ug/l	
189	ACID COMPOUNDS															
190	47	2-Chlorophenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
191	48	2,4-Dichlorophenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	4380	#####	#####	#####
192	49	2,4-Dimethylphenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	2020	365	#####	#####
193	50	4,6-Dinitro-o-Cresol	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
194	51	2,4-Dinitrophenol	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
195	52-53	Nitrophenols	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	230	150	#####	#####
196	54	4-Chloro-3-methylphenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	30	#####	#####	#####
197	55	Pentachlorophenol	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9.16	5.78	82	#####
198	56	Phenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	10200	2560	4600000	#####
199	57	2,4,6-Trichlorophenol	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	65	#####
201																
202	BASE/NEUTRAL COMPOUNDS															
203	58	Acenaphthene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1700	520	#####	#####
204	59	Acenaphthylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
205	60	Anthracene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
206	61	Benzidine	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	2500	#####	110000	#####
207	62	Benzol(a)anthracene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	5.40E-03	#####
208	63	Benzol(a)pyrene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
209	64	3,4-benzofluoranthene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
210	65	Benzo(g,h,i)perylene	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
211	66	Benzo(k)fluoranthene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
212	67	Bis(2-chloroethoxy)meth	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
213	68	Bis(2-chloroethyl) Ether	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14	#####
214	69	Bis(2-Chloroisopropyl) e	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	170E+05	#####
215	70	Bis(2-ethylhexyl)phthalate	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	59	#####
216	71	4-Bromophenyl phenyl e	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
217	72	Butylbenzyl phthalate	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
218	73	2-chloronaphthalene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1600	#####	#####	#####
219	74	4-chlorophenyl phenyl e	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
220	75	Chrysene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
221	76	Dibenzo(a,h)anthracene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
222	77-79	Dichlorobenzene(1,2	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	1120	763	#####	#####
223	80	3,3'-Dichlorobenzidine	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.77	#####
224	81	Diethyl Phthalate	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	1.20E+05	#####
225	82	Dimethyl phthalate	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	2.90E+06	#####
226	83	Di-n-Butyl phthalate	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
227	84	2,4-Dinitrotoluene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	330	230	91	#####
228	85	2,6-Dinitrotoluene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
229	86	Di-n-octyl phthalate	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
230	87	1,2-diphenylhydrazine	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	270	#####	#####	#####
231	88	Fluoranthene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	3980	#####	370	#####
232	89	Fluorene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14000	#####
233	90	Hexachlorobenzene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.0077	#####
234	91	Hexachlorobutadiene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	90	9.3	500	#####
235	92	Hexachlorocyclopentadi	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	7	5.2	1.74E+04	#####
236	93	Hexachloroethane	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	980	540	89	#####
237		Hexachlorocyclohexane	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	2	0.08	#####	#####
238	94	Indeno(1,2,3-cd)pyrene	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
239	95	Isophorone I	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	117000	#####	6000	#####
240	96	Naphthalene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	2300	620	#####	#####
241	97	Nitrobenzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	27000	#####	1800	#####
242	98	N-nitrosodimethylamine	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	81	#####
243	99	N-nitrosodi-n-propylamir	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
244	100	N-nitrosodiphenylamin	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	160	#####
245	101	Phenanthrene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
246	103	1,2,4-Trichlorobenzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
247																
248	PESTICIDES															
249	104	Aldrin	1	0.01	2.13	0	0	0.00	0.00	0.00	0.00	#####	3.00	#####	0.0014	#####
250	105	Alpha-BHC	1	0.05	2.13	0	0.0501	0.11	0.11	0.11	0.11	#####	2.00	0.08	0.0373	#####
251	106	Beta-BHC	1	0.05	2.13	0	0	0.00	0.00	0.00	0.00	#####	2.00	0.08	0.46	#####
252	107	Gamma-BHC	1	0.05	2.13	0	0.0642	0.14	0.14	0.14	0.14	#####	2.00	0.08	0.63	#####
253	108	Delta-BHC	1	0.05	2.13	0	0.0688	0.15	0.15	0.15	0.15	#####	2.00	0.08	#####	#####
254	109	Chlordane	1	0.2	2.13	0	0	0.00	0.00	0.00	0.00	#####	2.40	0.0043	0.005	#####
255	110	4,4'-DDT	1	0.02	2.13	0	0	0.00	0.00	0.00	0.00	#####	1.10	0.001	0.0059	#####
256	111	4,4'-DDE	1	0.1	2.13	0	0	0.00	0.00	0.00	0.00	#####	1.10	0.001	0.0059	#####
257	112	4,4'-DDD	1	0.1	2.13	0	0	0.00	0.00	0.00	0.00	#####	1.10	0.001	0.0084	#####





**Henderson, Katie**

---

**From:** Reiber, Loretta  
**Sent:** Tuesday, June 01, 2010 8:48 AM  
**To:** Henderson, Katie  
**Subject:** FW: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 4 of 4)

-----Original Message-----

**From:** Tillman.Michael@epamail.epa.gov [mailto:Tillman.Michael@epamail.epa.gov]  
**Sent:** Tuesday, August 04, 2009 9:43 AM  
**To:** bruce.fielding@la.gov  
**Cc:** Reiber, Loretta; Bailey, John; Baskin.Kilty@epamail.epa.gov  
**Subject:** Fw: GP Crossett, NPDES Permit No. AR0001210 - WQ screens (Part 4 of 4)

----- Forwarded by Michael Tillman/R6/USEPA/US on 08/04/2009 09:52 AM -----

**FW: GP Crossett, NPDES Permit No. AR0001210**



Reiber, Loretta

to: Michael Tillman

07/29/2009 02:42 PM

-----Original Message-----

**From:** Reiber, Loretta  
**Sent:** Wednesday, July 29, 2009 2:38 PM  
**To:** 'Tillman.Michael@epamail.epa.gov'  
**Subject:** GP Crossett, NPDES Permit No. AR0001210

Mike,

Attached are the Priority Pollutant Scan calculations and ADEQ's review. ADEQ is requesting a review of these documents in order to confirm compliance with LDEQ's procedures at SMS2.

If you have any questions, please feel free to contact me at (501) 682-0612 or by e-mail at [reiber@adeq.state.ar.us](mailto:reiber@adeq.state.ar.us).

Loretta Reiber, P.E.  
Engineer, NPDES Permits

A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
1					CALCULATIONS OF ARKANSAS WATER QUALITY-BASED EFFLUENT LIMITATIONS										
2					For an Arkansas River/Stream										
3							(Reserved)								
4	STEP 1:						GC								
5							Quachita River								
6	FACILITY														
7															
8															
9	Permittee						GP Crosssett								
10	NPDES Permit No.						AR0001210								
11	Outfall No. (s)						SMSZ								
12	Plant Effluent Flow (MGD)						52.40								
13	Plant Effluent Flow (cfs)						80.96								
14															
15	RECEIVING STREAM														
16															
17	Is this a large river? (see list at right)(enter "1" if yes, "0" if no, make entry as a number)						1								
18	Name of Receiving Stream:						Quachita River								
19	Waterbody Segment Code No						20								
20	Is this a lake or reservoir? (enter "1" if yes, "0" if no, make entry as a number)						0								
21	Second Enter 7Q10 in Cell H31						0								
22	(Reserved)						(Reserved)								
23	(Reserved)						(Reserved)								
24	(Reserved)						(Reserved)								
25	(Reserved)						(Reserved)								
26	(Reserved)						(Reserved)								
27	(Reserved)						(Reserved)								
28	(Reserved)						(Reserved)								
29	Ecoregion TSS (mg/l) (For Large River, See List to Right)						5.50								
30	Ecoregion Hardness (mg/l)						28.00								
31	Enter 7Q10 (cfs) as the Critical Flow (Reserved)						1200.00	(Reserved)							
32	Long Term Ave / Harmonic Mean Flow (cfs)						3600.00	(Reserved)							
33	Using Diffusers (Yes/No)						no								
34	pH (Avg)						7.01								
35	Percent (%) of Critical Flow for Chronic Criteria						0.25								
36	Percent (%) of Critical Flow for Acute Criteria						0.06								
37	Water Effect Ratio (WER)						1.00								
38	Ave Monthly Limit LTA Multiplier (Ref: page 103 TSD for WQ-Based Toxics Control)						1.55								
39	Max Daily Limit LTA Multiplier (Ref: page 103 TSD for WQ-Based Toxics Control)						3.11								

**Codes & TSS for Ecoregions and Large Rivers**

Arkansas (Fl. Smith to Dardanelle Dam) = 2.0 mg/l  
 Arkansas (Dardanelle Dam to Terry L&L) = 2.5 mg/l  
 Arkansas (Terry L&L to L&D No. 5) = 1.3 mg/l  
 Arkansas (L&D No. 5 to Mouth) = 3.0 mg/l  
 White (Above Beaver Lake) = 5.5 mg/l  
 White (From Bull Shoals to Black Riv) = 8.0 mg/l  
 White (From Black River to Mouth) = 18.5 mg/l  
 St. Francis River = 18.0 mg/l  
 Ouachita (Above Caddo River) = 2.0 mg/l  
 Ouachita (Below Caddo River) = 5.5 mg/l  
 Red River = 33.0 mg/l

**Total Hardness for:**  
 Arkansas River = 125 mg/l  
 Ouachita River = 28 mg/l  
 White River = 116 mg/l

Gulf Coastal = 31 mg/l  
 Ozark Highlands = 148 mg/l  
 Boston Mount = 25 mg/l

**Large Rivers**  
 Mississippi River, Arkansas River, Red River  
 White (Below confluence with Black River)  
 Ouachita (Below confluence with Little Miss. River)

For industrial and federal facility, use the highest monthly average flow for the past 24 months. For POTWs, use the design flow.

#VALUE! => No violation or Not Applicable  
 9999999.00 => No EPA/ADEQ Guideline



	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
40																
41	STEP 2:	INPUT AMBIENT AND EFFLUENT DATA														
42		CALCULATE IN-STREAM WASTE CONCENTRATIONS														
43																
44	DATA INPUT															
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70	Total Metals															
71		Dissolved Value in Stream	Kp	alpha (a)	Kp	C/Ct	C/Ct	Total Value			Kp	alpha (a)	Kp	C/Ct	Total Value	
72	Arsenic	480000	-0.73	138285.446	0.56799788	0.00	0.00	0.00			480000.00	-0.73	138285.45	0.5679979	0	
73	Cadmium	4000000	-1.13	582706.889	0.237818469	0.00	0.00	0.00			3520000.00	-0.92	733514.98	0.1986361	0	
74	Chromium(3)	3360000	-0.93	688338.365	0.208948818	0.00	0.00	0.00			2170000.00	-0.27	1369499.28	0.1172024	0	
75	Copper	1040000	-0.74	294554.016	0.381672529	0.00	0.00	0.00			2850000.00	-0.9	614495.12	0.2283249	0	
76	Lead	2800000	-0.8	715925.58	0.202527926	0.00	0.00	0.00			2040000.00	-0.53	826490.64	0.1803199	0	
77	Mercury	490000	-1.14	415321.613	0.30448177	0.00	0.00	0.00			1970000.00	-1.17	268066.09	0.4041443	0	
78	Nickel	1250000	-0.57	185433.992	0.495077211	0.00	0.00	0.00			2210000.00	-0.76	604946.03	0.2310962	0	
79	Zinc	3340000	-0.7	379014.766	0.324193117	0.00	0.00	0.00			3340000.00	-0.68	1047651.74	0.1478593	0	
80	Silver	2400000	-1.03	414607.994	0.30484608	0.00	0.00	0.00			2400000.00	-1.03	414607.99	0.3048461	0	
81																

\*Note: Use this section to convert lab concentrations shown as "Dissolved" to "Total"



A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
132	POLLUTANTS	Number of Data points	MQL ug/l	EPA Statistical Factor	Background Conc. ug/l	Effluent Conc. ug/l	Domestic Supply IWC ug/l	Acute Aquatic IWC ug/l	Chronic Aquatic IWC ug/l	Bioacc. IWC ug/l	Domestic Criteria ug/l	Arkansas Acute Criteria ug/l	Arkansas Chronic Aquatic ug/l	Arkansas Bioacc. ug/l	
133															
137	<b>METALS AND CYANIDE</b>														
138	1. Antimony Total	1	60	2.13	0	0	0.00	0.00	0.00	0.00	#####	9000.00	1600.00	4300	
139	2. Arsenic Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	50	633.81	334.51	1.4	
140	3. Barium Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	#####	130.00	5.30	4	
141	4. Cadmium Total	1	1	2.13	0	1.17	0.16	1.32	0.53	0.05	10	3.91	1.69	#####	
142	6. Chromium (Hex)	1	10	2.13	0	5.66	0.76	6.38	2.56	0.27	50	925.86	300.34	#####	
143	7. Chromium (Tri)	1	10	2.13	0	5.66	0.76	6.38	2.56	0.27	50	15.71	10.58	#####	
144	8. Copper Total	1	0.5	2.13	0	11.7	1.58	13.19	5.30	0.55	#####	13.44	10.02	#####	
145	9. Lead Total	1	0.5	2.13	0	3.13	0.42	3.53	1.42	0.15	50	77.87	3.03	#####	
146	10. Mercury Total	1	0.005	2.13	0	0.009	0.00	0.01	0.00	0.00	2	6.70	0.012	0.15	
147	12. Nickel Total	1	0.5	2.13	0	7.87	1.06	8.87	3.56	0.37	#####	973.88	108.16	4600	
148	13. Selenium Total	1	5	2.13	0	11.3	1.52	12.74	5.11	0.53	10	20.00	5.00	#####	
149	14. Silver Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	50	1.27	#####	6.3	
150	15. Thallium Total	1	0.5	2.13	0	0	0.00	0.00	0.00	0.00	#####	1400.00	#####	#####	
151	16. Zinc Total	1	20	2.13	0	251	33.79	282.97	113.62	11.76	#####	120.05	109.63	#####	
152	129. Phenols Total	1	5	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
153	17. Cyanide Total	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	22.36	5.2	220000	
155	<b>DIOXIN</b>														
157	18. 2,3,7,8-TCDD	1	0.00001	2.13	0	0	0.00	0.00	0.00	0.00	#####	0.01	1.00E+07	1.00E-06	
159	<b>VOLATILE COMPOUNDS</b>														
160	18. Acrolein	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	68	21	780	
161	20. Acrylonitrile	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	7550	2600	6.6	
162	21. Benzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	5300	9999999	710	
163	22. Bromoform	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	3600	
164	23. Carbon Tetrach	1	2	2.13	0	0	0.00	0.00	0.00	0.00	5	35200	#####	44	
165	24. Chlorobenzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	250	50	2.10E+04	
166	25. Chlorobromomethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	340	
167	26. Chloroethane	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	1.00E+07	
168	27. 2-Chloroethylvinyl ether	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999	#####	1.00E+07	
169	28. Chloroform	1	10	2.13	0	0	0.00	0.00	0.00	0.00	7	28900	1240	4700	
170	29. Dichlorobromomethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	220	
171	30. 1,1-Dichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	11800	20000	990	
172	31. 1,2-Dichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	11600	#####	32	
173	32. 1,1-Dichloroethylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	23000	5700	#####	
174	33. 1,2-Dichloropropane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	6060	244	1700	
175	34. 1,3-Dichloropropylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	32000	#####	29000	
176	35. Ethylbenzene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
177	37. Methyl Chloride	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
178	36. Methyl bromide	1	50	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	4000	
179	38. Methylene chloride	1	20	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	16000	
180	39. 1,1,2,2-Tetrachloroetha	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9320	2400	110	
181	40. Tetrachloroethylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	5280	840	88.5	
182	41. Toluene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	17500	#####	2.00E+05	
183	42. 1,2-trans-dichloroethyler	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	
184	44. 1,1,2-Trichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	#####	18000	9400	420	
185	43. 1,1,1-Trichloroethane	1	10	2.13	0	0	0.00	0.00	0.00	0.00	200	18000	#####	#####	
186	45. Trichloroethylene	1	10	2.13	0	0	0.00	0.00	0.00	0.00	5	45000	21900	810	
187	46. Vinyl Chloride	1	10	2.13	0	0	0.00	0.00	0.00	0.00	2	9999999.00	#####	5250	

TOX	A	B	C	D	E	F	G	H	I	J	K	L	M	N	O	P
						Ambient Background Conc.	Effluent Conc.	Domestic Supply IWC	Acute Aquatic IWC	Chronic Aquatic IWC	Human Health IWC	Domestic Criteria ug/l	Acute Aquatic Criteria ug/l	Chronic Aquatic Criteria ug/l	Human Health Criteria ug/l	
189	ACID COMPOUNDS															
190	47. 2-Chlorophenol			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
191	48. 2,4-Dichlorophenol			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	4380	#####	#####	#####
192	49. 2,4-Dimethylphenol			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	2020	365	#####	#####
193	50. 4,6-Dinitro-o-Cresol			1	50	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
194	51. 2,4-Dinitrophenol			1	50	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	765	#####
195	52. 4,6-Dinitrophenol			1	50	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14000	#####
196	54. 4-Chloro-3-methylphenol			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	230	150	#####	#####
197	55. Pentachlorophenol			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	30	#####	#####	#####
198	56. Phenol			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9.16	5.78	82	#####
199	57. 2,4,6-Trichlorophenol			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	10200	2560	4600000	#####
200																
201	BASE/NEUTRAL COMPOUNDS															
202	58. Acenaphthene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	1700	520	#####	#####
203	59. Acenaphthylene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
204	60. Anthracene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
205	61. Benzidine			1	50	2.13	0	0.00	0.00	0.00	0.00	#####	2500	#####	110000	#####
206	62. Benz(a)anthracene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	5.40E-03	#####
207	63. Benzo(a)pyrene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
208	64. 3,4-Benzofluoranthene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
209	65. Benzo(g,h,i)perylene			1	20	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
210	66. Benzo(k)fluoranthene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
211	67. Bis(2-chloroethoxy)meth			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
212	68. Bis(2-chloroethyl) ether			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
213	69. Bis(2-chloroisopropyl) e			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14	#####
214	70. Bis(2-ethylhexyl)phthal			1	10	2.13	0	0.39	3.24	1.30	0.13	#####	9999999.00	#####	1.70E+05	#####
215	71. 4-Bromophenyl phenyl e			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	59	#####
216	72. Butylbenzyl phthalate			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
217	73. 2-chloronaphthalene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	1600	763	#####	#####
218	74. 4-chlorophenyl phenyl e			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
219	75. Chrysene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
220	76. Dibenz(a,h)anthracene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
221	77. 7,9-Dichlorobenzene(1,2			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	1120	#####	#####	#####
222	80. 3,3'-Dichlorobenzidine			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	2600	#####
223	81. Diethyl phthalate			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.77	#####
224	82. Dimethyl phthalate			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	1.20E+05	#####
225	83. Di-n-Butyl phthalate			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	2.90E+06	#####
226	84. 2,4-Dinitrotoluene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	1.20E+04	#####
227	85. 2,6-Dinitrotoluene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	91	#####
228	86. Di-n-octyl phthalate			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
229	87. 1,2-diphenylhydrazine			1	20	2.13	0	0.00	0.00	0.00	0.00	#####	270	#####	5.4	#####
230	88. Fluoranthene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	3980	#####	370	#####
231	89. Fluorene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	14000	#####
232	90. Hexachlorobutadiene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.0077	#####
233	91. Hexachlorocyclopentadi			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	90	9.3	500	#####
234	92. Hexachlorocyclohexane			1	20	2.13	0	0.00	0.00	0.00	0.00	#####	7	5.2	1.74E+04	#####
235	93. Indene(1,2,3-c)pyrene			1	5	2.13	0	0.00	0.00	0.00	0.00	#####	2	0.08	#####	#####
236	94. Isophorone			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	0.31	#####
237	95. Naphthalene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	117000	#####	6000	#####
238	96. Nitrobenzene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	2300	620	#####	#####
239	97. Nitrosodimethylamine			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	27000	#####	1900	#####
240	98. N-nitrosodiphenylamine			1	50	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	81	#####
241	99. N-nitrosodipropylamine			1	20	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
242	100. N-nitrosodiphenylamin			1	20	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	160	#####
243	101. Phenanthrene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
244	103. 1,2,4-trichlorobenzene			1	10	2.13	0	0.00	0.00	0.00	0.00	#####	9999999.00	#####	#####	#####
245																
246	PESTICIDES															
247	104. Aldrin			1	0.01	2.13	0	0.00	0.00	0.00	0.00	#####	3.00	#####	0.0014	#####
248	105. Alpha-BHC			1	0.05	2.13	0	0.00	0.00	0.00	0.00	#####	2.00	0.08	0.0373	#####
249	106. Beta-BHC			1	0.05	2.13	0	0.00	0.00	0.00	0.00	#####	2.00	0.08	0.46	#####
250	107. Gamma-BHC			1	0.05	2.13	0	0.01	0.11	0.04	0.00	#####	2.00	0.08	0.63	#####
251	108. Delta-BHC			1	0.05	2.13	0	0.01	0.05	0.02	0.00	#####	2.00	0.08	#####	#####
252	109. Chlordane			1	0.2	2.13	0	0.00	0.00	0.00	0.00	#####	2.40	0.0043	0.005	#####
253	110. 4,4'-DDT			1	0.02	2.13	0	0.00	0.00	0.00	0.00	#####	1.10	0.001	0.0059	#####
254	111. 4,4'-DDE			1	0.1	2.13	0	0.00	0.00	0.00	0.00	#####	1.10	0.001	0.0059	#####
255	112. 4,4'-DDD			1	0.1	2.13	0	0.00	0.00	0.00	0.00	#####	1.10	0.001	0.0084	#####



