

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

Figure 4-6 - Observed Vapor Pressure of 48.8% Aqueous Acrylamide

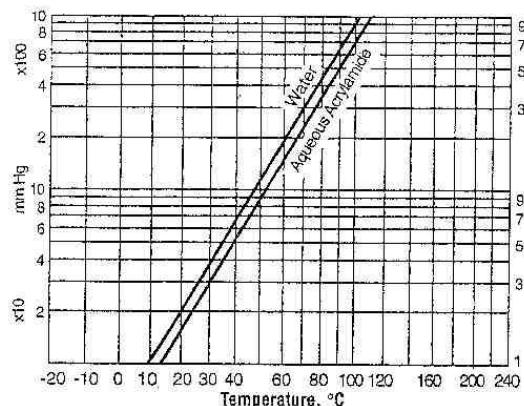


Table 4-4 - Vapor Phase Concentrations of AAm in PPM (100% Solid and 50% Aqueous AAm)

Temperature °C	Vapor Phase Concentration AAm over 100% Solid AAm	Vapor Phase Concentration AAm over 50% Aqueous AAm
10	6.2	1.3
20	16.6	3.5
25	26	5.6
30	41	9
37	75	17
80	1700	430

Table 4-5 - Vapor Phase Concentrations of AAm in PPM of Different AAm Solutions

°C	mm Hg
23	18.05
31	29.79
38	43.36
53	92.72
70	209.49

Temperature °C	Weight Percent AAm in Aqueous Solutions				
	10	18	25	35	50
10	0.35	0.57	0.74	0.94	1.3
20	0.94	1.6	2.0	2.6	3.6
25	1.5	2.5	3.2	4.1	5.7
30	2.4	4.0	5.1	6.5	9.1
37	4.4	7.3	9.5	12	17
50	13	22	27	38	54
70	56	94	114	158	225
80	108	180	219	303	431
90	198	332	405	560	769
100	353	591	721	997	1418

Saturated Vapor Phase Concentrations of AAm in PPM as a Function of Temperature for Different Aqueous Concentrations (Including Pure Crystal)

Experiments have been conducted to determine the vapor phase concentration in ppm of pure crystal, as well as for different concentrations of crystal acrylamide monomer in aqueous solutions as a function of temperature. It is important to recognize that these values were obtained from vapor-liquid equilibrium experiments in a closed container. Any departure from equilibrium, such as produced in an open vessel, or from continuous air-sparging throughout the monomer liquid in an acrylamide monomer storage tank, will result in much less than the reported equilibrium vapor phase concentration.

Tables 4-4 and 4-5 compare the vapor phase concentrations of solid acrylamide monomer and that of various weight % acrylamide monomer solutions as a function of temperature.

For PAM Reactor, actual conditions are ~5 wt% and 95°F (35°C)

use data @ 10wt%, 37°C

$$y_i = 4.4 \text{ ppm} / 10^6 \\ = 4.4 \times 10^{-6} \text{ mol/mole}$$

$$\text{partial press} = \gamma P = y_i P$$

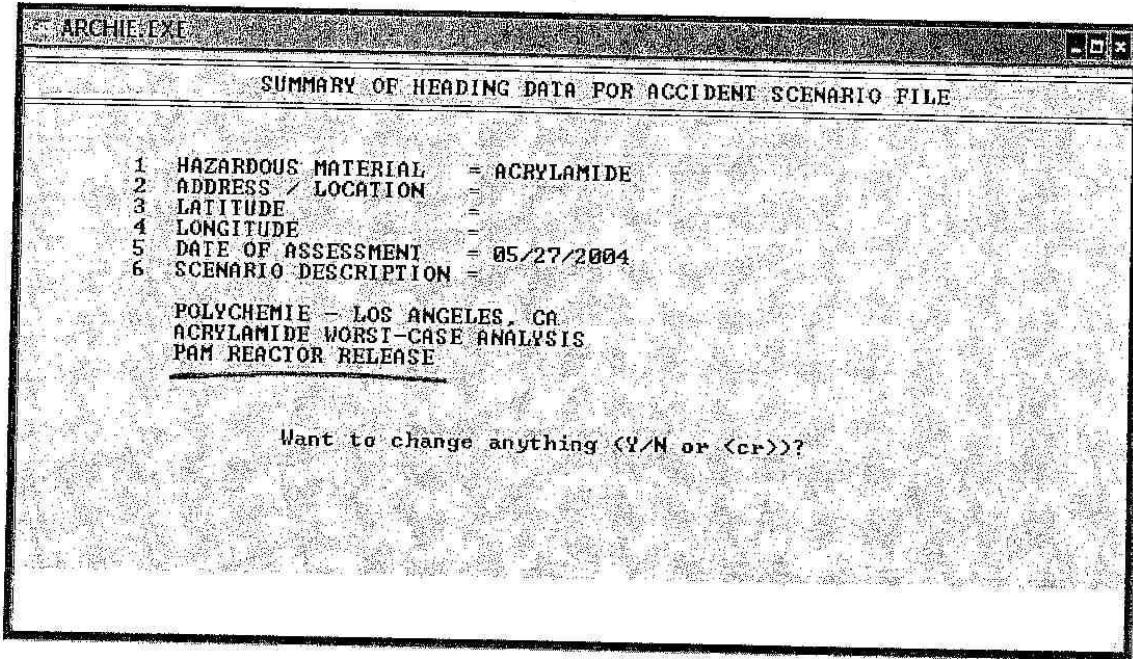
$$\gamma P = 4.4 \times 10^{-6} \times 760 \text{ mmHg} \\ = 0.00334 \text{ mmHg}$$

for ADAM-Quat RXR, actual cond. are ~5 wt%, 131°F (55°C)

use data @ 10wt% and 70°C

$$y_i = 56 \text{ ppm} / 10^6$$

$$P = \frac{56}{10^6} \times 760 \text{ mmHg} \\ = 0.0426 \text{ mmHg}$$



V

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

PAM Reactor Release

ARCHIE-EXE

CURRENT PARAMETER VALUES FOR LIQUID POOL EVAPORATION RATE ESTIMATION		
1	MOLECULAR WEIGHT	= 70.07
2	NORMAL BOILING POINT	= 212 degrees F
3	LIQUID SPECIFIC GRAVITY	= 1.04
4	LIQ TEMP IN TANK	= 112 degrees F
5	AMBIENT TEMPERATURE	= 112 degrees F
6	LIQ VAP PRES AT AMB TEMP	= .00334 mm Hg
7	EVAPORATING POOL AREA	= 2819 sq ft
8	WEIGHT SPILLED LIQUID	= 94800 lbs
9	ATMOS STABILITY CLASS	= F
10	WIND VELOCITY	= 3.4 mph

MODEL RESULTS:

Vapor evolution rate = .002 lbs/min
 Vapor evolution duration = 5.471698E+07 minutes

Want to rerun the model with different input values (Y/N or <cr>)?

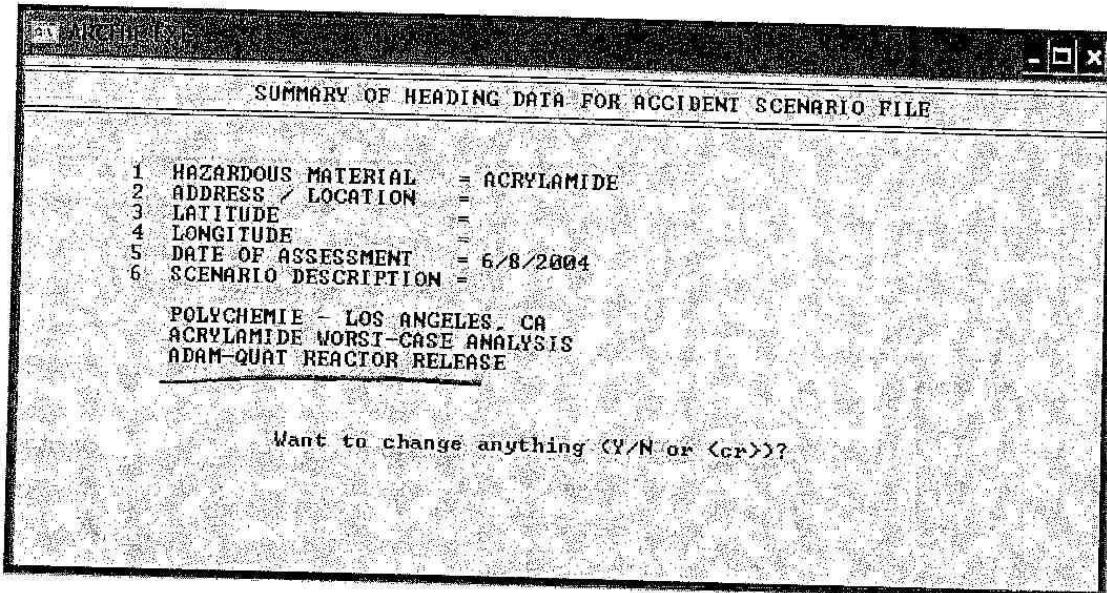
ARCHIE-EXE

CURRENT PARAMETER VALUES FOR TOXIC GAS OR VAPOR HAZARD EVALUATION		
1	MOLECULAR WEIGHT	= 70.07
2	TOXIC VAPOR LIMIT	= 40.91 ppm
3	VAPOR/GAS DISCHARGE HEIGHT	= 0 feet
4	ATMOSPHERIC STABILITY CLASS	= F
5	WIND VELOCITY AT SURFACE	= 3.4 mph
6	AMBIENT AIR TEMPERATURE	= 112 deg F
7	VAPOR/GAS EMISSION RATE	= .002 lb/min
8	DURATION OF EMISSION	= 10 minutes

MODEL RESULTS:

Downwind toxic hazard distance
 on ground (min. computable) < 33 feet
 ← 0.1 miles

Tables of concentrations and times vs distance not available!
 Want to rerun the model with different input values (Y/N or <cr>)?



APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

ADAM-Qut Reactor Release

CURRENT PARAMETER VALUES FOR LIQUID POOL EVAPORATION RATE ESTIMATION		
1	MOLECULAR WEIGHT	= 70.07
2	NORMAL BOILING POINT	= 212 degrees F
3	LIQUID SPECIFIC GRAVITY	= 1.04
4	LIQ TEMP IN TANK	= 131 degrees F
5	AMBIENT TEMPERATURE	= 112 degrees F
6	LIQ VAP PRES AT AMB TEMP	= .0426 mm Hg
7	EVAPORATING POOL AREA	= 2819 sq ft
8	WEIGHT SPILLED LIQUID	= 17525 lbs
9	ATMOS STABILITY CLASS	= F
10	WIND VELOCITY	= 3.4 mph
MODEL RESULTS:		
Vapor evolution rate	= .028	lbs/min
Vapor evolution duration	= 640784.8	minutes
Want to rerun the model with different input values (Y/N or <cr>)?		

CURRENT PARAMETER VALUES FOR TOXIC GAS OR VAPOR HAZARD EVALUATION		
1	MOLECULAR WEIGHT	= 70.07
2	TOXIC VAPOR LIMIT	= 40.91 ppm
3	VAPOR/GAS DISCHARGE HEIGHT	= 0 feet
4	ATMOSPHERIC STABILITY CLASS	= F
5	WIND VELOCITY AT SURFACE	= 3.4 mph
6	AMBIENT AIR TEMPERATURE	= 112 deg F
7	VAPOR/GAS EMISSION RATE	= .028 lb/min
8	DURATION OF EMISSION	= 10 minutes
MODEL RESULTS:		
Downwind toxic hazard distance at groundlevel	= 67	feet
<i><0.1 mile</i>		
*** Press ENTER to View Various Tables ***		

ADAM-Qxt Reactor Release

Downwind Distance (feet)	Ground Level Concentration (ppm)	Source Height Concentration (ppm)	Initial evacuation Zone Width (feet)
32	.01	190	54
35	.01	163	56
37	.01	140	57
40	.01	122	59
42	.01	108	60
45	.01	95.2	61
47	.01	84.9	63
50	.01	76.3	65
52	.01	68.9	66
55	.02	62.6	68
57	.02	57	70
60	.02	52.2	72
62	.02	48	74
65	.02	44.3	75
67	.02	40.91	76

Usually safe for 1-hour release. Longer releases or sudden wind shifts may require a larger width or different direction for the evacuation zone.
 ***** Press ENTER to Continue *****

Downwind Distance (feet)	Contaminant Arrival Time at Downwind Location (minutes)	Contaminant Departure Time at Downwind Location (minutes)
32	.01	10.2
35	.01	10.3
37	.01	10.3
40	.01	10.3
42	.01	10.3
45	.01	10.4
47	.01	10.4
50	.01	10.4
52	.01	10.4
55	.02	10.4
57	.02	10.4
60	.02	10.4
62	.02	10.4
65	.02	10.4
67	.02	10.4

CAUTION: See guide for assumptions used in estimating the scenario.
 Want to rerun the model with different input values? (CN or KO)

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

WORST-CASE SCENARIO FOR TOXIC LIQUIDS
 Polychemic Inc. - Los Angeles, CA
 GESSI Project No. 02512.01

Toxic Liquid	Maximum Quantity Released ¹ QS	Area of Dike ² (ft ²)	Density Factor ³ of Liquid DF	Maximum Area of Pool ⁴ (ft ²)	Liquid Factor Ambient ⁵ LFA	Release Rate ⁶ QR (lb/min)	Duration of Release ⁷ t (min)	Maximum Duration of Release ⁸ (min)	Toxic Endpoint ⁹ (mg/l)	Release/Toxic Endpoint Table	Reference Table to Use	Distance to Endpoint ¹⁰ (miles)
Acrylamide, density, lb/gal: 8.68	20,000 gals 173,600 lbs, 50% 86,800 lbs, 100%	NA	0.47	81,019	4.32E-06	0.49	354,532	10	0.11	4.5	3	0.1 (Shortest reportable distance)
Acrylamide, density, lb/gal: 8.68	781,200 lbs, 37% 390,600 lbs, 100%	RAILCAR RELEASE										
Acrylamide, density, lb/gal: 8.68	30,000 gals 260,400 lbs, 50% 130,200 lbs, 100%	TANK (proposed) RELEASE	0.47	2819	4.32E-06	0.02	15,284,022	10	0.11	0.15	3	0.1 (Shortest reportable distance)

NOTES:

- 1 - Calculations are based on U.S. EPA RMP Guidance for Offsite Consequence Analysis provided by the EPA dated April 15, 1999.
- 2 - Maximum quantity released is equal to the capacity of the largest vessel handling the liquid.
- 3 - For the release rate, use the minimum of the dike area (less tank area) or the area calculated using the density factor (see Note 4).
- 4 - Density Factor of Liquid, $DF = 1 / 0.023 * \text{density of substance} / (\text{ft}^3) = 1 / 0.023 * (8.68 \text{ lb/gal} * 0.74805 \text{ gal/ft}^3) [Eq. D-4]$
- 5 - Maximum area of the liquid pool, $A (\text{ft}^2) = \text{Quantity Released} \cdot QS(\text{lb}) * \text{Density Factor of the Liquid (DF)} [Eq. 3-6]$
- 6 - Liquid Factor Ambient (LFA) = $0.254 * MW (71.08) za * VP (0.02166 \text{ mmHg}) / 82.05*298 [Eq. D-2]$
- 7 - Release Rate (no passive mitigation), $QR = QS * 1.4 * LFA * DF [Eq. 3-7]$
- 8 - Duration of Release, $t = \text{Quantity Released}, QS (\text{lb}) / \text{Release Rate}, QR (\text{lb/min}) [Eq. 3-5]$
- 9 - Maximum Duration of the release is from the EPA RMP Guidance for OCA. The maximum duration for water solutions of toxic substances is 10 minutes. [Sect. 3.3]
- 10 - Toxic Endpoint from Appendix B of California Accidental Release Prevention (CalARP) Program. Administering Agency Guidance, California Governor's Office of Emergency Services (OES), August 28, 2003.
- 11 - Distance to endpoint is calculated using Reference Table.

By: NSB 05/27/2004
 Checked: KSD 05/28/2004

Polychemie Inc. – Facility Expansion Project



CALCULATION SHEET

Calc. No.

TS-2

Prepared By: Natasha Brash Date: 5/27/2004 Reviewed By: Karen Dorman Date: 5/27/2004

Section I. General Information

Project: Polychemie, Inc. - Los Angeles, CA
2004 RMP Update Project No.: 02512.01

Subject: RMP Alternate Case, Toxic Substance Release - Acrylamide

Section II. Scenario

- A. 50% acrylamide is released from a ruptured unloading hose on a railcar during delivery. Break occurs before pump (Option 1) or after pump (Option 2). Compare quantity released from both options and use larger amount in subsequent analysis. Response time to stop release is 3 minutes.

Section III. Calculations

- A. The OCA document (Ref. A) does not directly address releases for flammable liquids from a pressurized storage vessel. Therefore, the equation used for the alternate analysis for liquids in Appendix D of the OCA document will be used.

Option 1: Per OCA document, Equation D-13:

$$\text{Discharge rate, } m \text{ (kgs/sec)} = A_h C_d (\rho_l [2g p_i (H_L - H_h) + 2(P_0 - P_a)])^{0.5}$$

A_h = opening area (m^2) = assume hose = 3 in = 0.0762 m = 0.005 m^2
 C_d = discharge coefficient (unitless) = 0.62 From ARCHIE memo (Ref. B)
 g = gravitational constant = 9.8 m/s^2
 ρ_l = liquid density (kg/m^3) = 1000 kg/m^3 x 1.04 = 1040 kg/m^3
 P_0 = storage pressure = 25 psig = 39.7 psia = 273,693 Pa
 P_a = ambient pressure = 14.7 psia = 101,325 Pa
 H_L = liquid height above bottom of container = assume top of railcar = 10 ft = 3.048 m
 H_h = height of opening = 0 m (bottom loading)
 m = 58.16 kg/sec = 3490 kg/min = 7,693 lbs/min

Assume response time, t = 3 minutes

Quantity release, $QS = QF_L$ (or m) $\times t$ = 23,080 pounds Eq 3-5 of OCA Guidance

- B. Option 2: Approximate quantity released as amount being pumped.

Liquid being pumped at 40-60 gpm.
Use 60 gpm and increase by 1.25 to account for increase in inflow due to loss of head.

$$\text{Flow} = 60 \text{ gpm} \times 1.25 = 75 \text{ gpm}$$
$$QR = 75 \text{ gpm} \times 3 \text{ min} = 225 \text{ gals} \times 8.68 \text{ lb/gal} = 1953 \text{ lbs}$$

Section IV. Results

- A. Evaluate Option 1 since quantity released before pump is greater.

Section V. References

- A. Risk Management Program Guidance for Offsite Consequence Analysis. U.S. Environmental Protection Agency, Office of Solid Waste and Emergency Response, (EPA 550-B-99-009), 04/15/1999.
- B. "Correction of Program Error -- New Version of MODCON.EXE" memo. Memo issued as an update to the ARCHIE model referenced in The Handbook of Chemical Hazard Analysis Procedures, Department of Transportation, the Federal Emergency Management Agency, and the Environmental Protection Agency, (http://hazmat.dot.gov/risk_tools.htm), 1989.

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

ALTERNATE SCENARIO FOR TOXIC LIQUIDS
Polychemic Inc. - Los Angeles, CA
GESI Project No. 02512.01

Toxic Liquid	Quantity Released ¹ QS	Density Factor ² of Liquid DF	Liquid Factor ³ Ambient LFA	Release Rate ⁴ QR (lb/min)	Duration of Release ⁵ (min)	Maximum Duration of Release ⁶ (min)	Toxic Endpoint ⁷ (mg/l)	Release Rate/Toxic Endpoint ⁷	Reference Table to Use	Reference Table	Distance to Endpoint ⁸ (miles)
Acrylamide 50% density = 8.68 lb/gal	23,080 lbs, 50% 11,540 lbs, 100%	0.47	4.32E-06	0.11	206,811	10	0.11	1.0	Buoyant - Urban - 10 minute release	16	0.1 (Shortest reportable distance)

NOTES:

- Calculations are based on U. S. EPA RMP Guidance for Offsite Consequence Analysis provided by the EPA dated April 15, 1999.
- 1 - Quantity Released is assumed to equal spill release amt - see Calculation No. TS-2
- 2 - Density Factor of Liquid, $DF = 1 / 0.033 * \text{density of substance}(\text{lb/ft}^3) = 1 / 0.033 * (8.68 \text{ lb/gal} * 0.74805 \text{ gal/ft}^3)$ [Eq. D-4]
- 3 - Liquid Factor Ambient (LFA) = $0.284 * \text{MW} (71.08)^{0.25} * \text{VP} (0.02166 \text{ mmHg}) / 82.05^{*298}$ [Eq. D-2]
- 4 - Unmitigated Release Rate, QR = QS * 2.4 * LFA * DF [Eq. 7-10]
- 5 - Duration of Release, t = Quantity Released(QS lb) / Release Rate, QR (lb/min) [Eq. 3-5]
- 6 - Maximum Duration of the release is from the EPA RMP Guidance for OCA. The maximum duration for water solutions of toxic substances is 10 minutes. [Sect. 3.3]
- 7 - Toxic Endpoint from Appendix B of California Accidental Release Prevention, (CalARP) Program: Administering Agency Guidance, California Governor's Office of Emergency Services (OES), August 28, 2003.
- 8 - Distance to endpoint is calculated using Reference Table.

By: NSB 05/27/2004
Checked: KSD 05/27/2004

Vapor Phase Concentrations of Acrylamide (ppm_v) at Various Temperatures
50% Aqueous Solution

Temperature		Vapor Concentration	Equivalent Vapor Pressure	Equivalent Vapor Pressure
(°C)	(°F)	(ppm _v)	(psia)	(mmHg)
10	50	1.3	0.000096	0.00494
20	68	3.6	0.000265	0.01368
25	77	5.7	0.000419	0.02166
30	86	9.1	0.000669	0.03458
37	98.6	17	0.001249	0.06460
50	122	54	0.003968	0.20520
70	158	225	0.016533	0.85500
80	176	431	0.031670	1.63780
90	194	769	0.056506	2.92220
100	212	1418	0.104195	5.38840

Vapor concentration - temperature data provided by Dow Chemical

Vapor pressure estimated as follows (p_a):

$$\text{ppm}_v / 1 \times 10^6 = \text{vapor mole fraction } y_a$$

$$x_a = \text{liquid mole fraction}$$

$$p_a (\text{psia}) = y_a P / x_a, \text{ where } P = 14.696 \text{ psia}$$

Molecular weight acrylamide =

70.07 lb/lb mole

By:

RDH

Molecular weight water =

18 lb/lb mole

Checked:

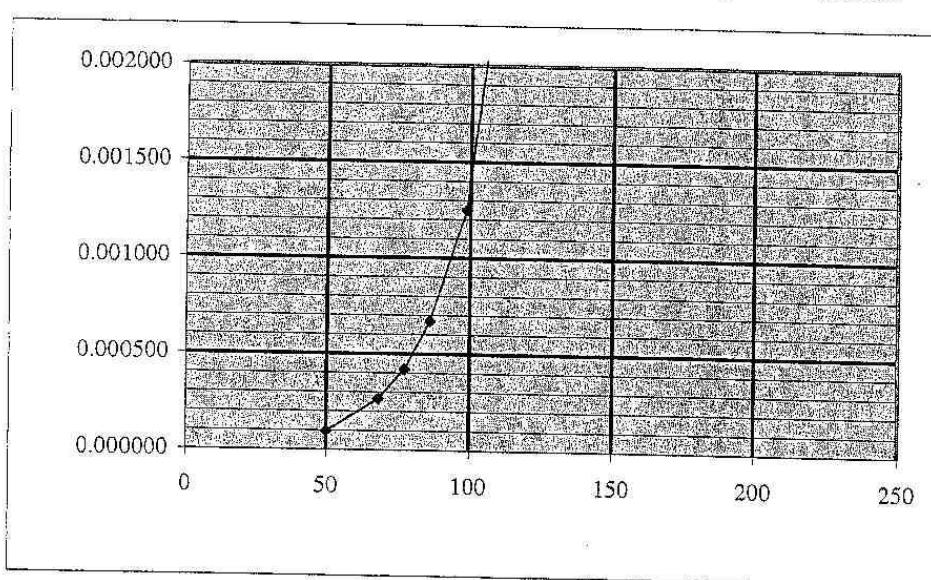
KSD

For a 50% aqueous solution, x_a =

0.2044

Date:

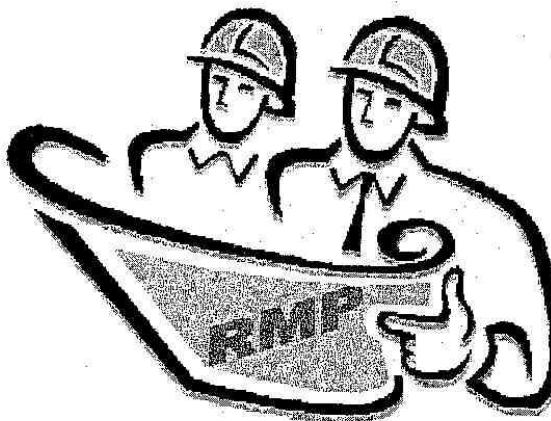
7/1/1998





California Accidental Release Prevention (CalARP) Program

Administering Agency Guidance



August 28, 2003

Appendix B
CalARP Program Toxic Endpoint Table

The following Toxic Endpoint (TE) Table should be used for all toxic substances listed in Section 2770.5, Table 1 and Table 3, of the CalARP Program. Where USEPA provided a TE for the FedARP Program, that TE is listed in the Table below. All other TEs were provided by the Office of Environmental Health Hazard Assessment (OEHHA), using preexisting toxicity values.

Table of Toxic Endpoints
[to be used as described in Section 2750.2 of the CalARP Program regulations]

Chemical Name	CAS Number	Toxic Endpoint (TE)		Basis for TE
		TE in (gm/m ³ or mg/l)	TE in (ppm) ¹	
Acetone Cyanohydrin	75-86-5	0.012	3	USEPA LOC ²
Acetone Thiosemicarbazide	1752-30-3	0.1		USEPA LOC ²
Acrolein [2-Propenal]	107-02-8	0.0011	0.5	USEPA ARP Program ³
Acrylamide	79-06-1	0.11		USEPA LOC ²
Acrylonitrile [2-Propenenitrile]	107-13-1	0.076	35	USEPA ARP Program ³
Acrylyl Chloride [2-Propenyl Chloride]	814-68-6	0.00090	0.2	USEPA ARP Program ³
Aldicarb	116-06-3	0.0003		USEPA LOC ²
Aldrin	309-00-2	0.002		IDLH95/10 ²
Allyl Alcohol [2-Propen-1-ol]	107-18-6	0.036	15	USEPA ARP Program ³
Allylamine [2-Propen-1-amine]	107-11-9	0.0032	1	USEPA ARP Program ³
Aluminum Phosphide	20859-73-8	0.02		USEPA LOC ²
Aminopterin	54-62-6	0.025		USEPA LOC ²
Amiton Oxalate	3734-97-2	0.003		USEPA LOC ²
Ammonia (anhydrous) or (aqueous), or Ammonium Hydroxide	7664-41-7	0.14	200	USEPA ARP Program ³
Aniline	62-53-3	0.038	10	USEPA LOC ²
Antimycin A	1397-94-0	0.0018		USEPA LOC ²
ANTU	86-88-4	0.01		USEPA LOC ²
Arsenic Pentoxide	1303-28-2	0.0005 as As		IDLH95/10 ²
Arsenous Oxide	1327-53-3	0.0005 as As		IDLH95/10 ²
Arsenous Trichloride	7784-34-1	0.010	1	USEPA ARP Program ³
Arsine	7784-42-1	0.0019	0.6	USEPA ARP Program ³
Azinophos-Ethyl	2642-71-9	0.0039		USEPA LOC ²
Azinophos-Methyl	86-50-0	0.001		IDLH95/10 ²
Benzene, 1-(Chloromethyl)-4-Nitro-	100-14-1	0.028		USEPA LOC ²
Benzeneearsonic Acid	98-05-5	0.00027		USEPA LOC ²
Benzimidazole, 4,5-Dichloro-2-(Trifluoromethyl)-	3615-21-2	0.013		USEPA LOC ²
Benzotrichloride	98-07-7	0.0007	0.1	USEPA LOC ²
Bicyclo[2.2.1] Heptane-2-Carbonitrile, 5-Chloro- 6-(((Methylamino) Carbonyl)Oxy)Imino)-, (1s-(1-alpha, 2-beta, 4-alpha, 5-alpha, 6E))-	15271-41-7	0.019		USEPA LOC ²
Bis(Chloromethyl) Ketone	534-07-6	0.00027		USEPA LOC ²
Bitoscanate	4044-65-9	0.02		USEPA LOC ²
Boron Trichloride [Borane, Trichloro-]	10294-34-5	0.010	2	USEPA ARP Program ³
Boron Trifluoride [Borane, Trifluoro-]	7637-07-2	0.028	10	USEPA ARP Program ³
Boron Trifluoride compound w/ Methyl Ether (1:1) [Boron, Trifluoro [oxybis[methane]]]-, T-4-	353-42-4	0.023	5	USEPA ARP Program ³
Bromadiolone	28772-56-7	0.001		USEPA LOC ²

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

Page 1 of 1

KCQT Extremes
Highest Maximum Temp
JAN 1 - DEC 31 1921 - 2002

INDIVIDUAL DAYS degrees	date	365/366 DAY PERIOD	
		degrees	year
1. 112	6-26-1990	1.	77.7 1981
2. 110	9-04-1988	2.	77.6 1984
3. 110	9-01-1955	3.	77.5 1986
4. 109	6-27-1990	4.	77.3 1997
5. 109	9-26-1963	5.	77.0 1992
6. 108	10-04-1987	6.	76.9 1980
7. 108	10-03-1987	7.	76.9 1994
8. 108	9-02-1955	8.	76.8 1990
9. 107	10-10-1991	9.	76.8 1995
10. 107	7-01-1985	10.	76.7 1959
11. 107	9-25-1978	11.	76.7 1983
12. 107	9-25-1963	12.	76.7 1996
13. 107	9-20-1939	13.	76.6 1958
14. 106	4-06-1989	14.	76.3 1972
15. 106	9-24-1978	15.	76.3 1985
16. 106	6-20-1973	16.	76.1 1989
17. 106	9-13-1971	17.	76.1 1993
18. 106	9-28-1963	18.	75.6 1967
19. 106	9-27-1963	19.	75.5 1987
20. 106	9-21-1939	20.	75.4 1966

Total Years = 82
Missing days = 336

Los Angeles
Civic Center

Polychemie Inc. – Facility Expansion Project

*Review of Worst Case and Alternative Release Scenarios
Polychemie Inc. – Los Angeles, California
GESI Project No. 02512.01*

June 2004

APPENDIX B

Impacted Population and Environmental Receptor Data

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS



CALCULATION SHEET

Calc. No.

Pop-1

Prepared By: Natasha Brash Date: 6/3/2004 Reviewed By: Karen Dorman Date: 6/3/2004

Section I. General Information

Project: Polychemie Inc. - Los Angeles, CA
2004 RMP Update

Project No.: 02512.01

Subject: Impacted Residential Population Estimates

Section II. Background

- A. From off-site consequence analysis, impacted areas are at 0.1, 0.2, and 0.32 mile radii. Total population within these impacted radii are estimated using aerial photographs, U.S. Census data, and general knowledge of the area. Aerial photographs and U.S. Census data showing total residential population per block are attached.

Section III. Estimates

- 0.1 mile radius: From general knowledge of the area, only industries are located within a 0.1 mile radius. Therefore, residential population is estimated to be 0.
- 0.2 mile radius: Beyond the immediate industrial area, there are residential areas to the southwest and north at the 0.2 mile radius. From the U.S. Census data showing total residential population per block, there are approximately 2 blocks with a population of 100-250 persons/block. Using 250 as a conservative estimate, total residential population is equal to $2 \times 250 = 500$ persons.
- 0.32 mile radius: At 0.32 mile radius, the residential population expands to the northeast in addition to the southwest and north. Using U.S. Census data showing total residential population per block, total residential population is estimated as follows:

$$\begin{aligned} 1 \text{ block} \times 50 \text{ persons} &= 50 \\ 2\frac{1}{2} \text{ blocks} \times 100 \text{ persons} &= 250 \\ 3\frac{1}{2} \text{ blocks} \times 250 \text{ persons} &= 875 \\ 1\frac{1}{2} \text{ blocks} \times >250 \text{ persons} & \\ (\text{Use } 3 \text{ blocks} \times 250 \text{ as conservative estimate}) \\ 3 \text{ blocks} \times 250 \text{ persons} &= 750 \\ \text{Total residential population} &= 1925 \end{aligned}$$

Section IV. Results

Radius (miles)	Total Population
0.1	0
0.2	500
0.32	1925



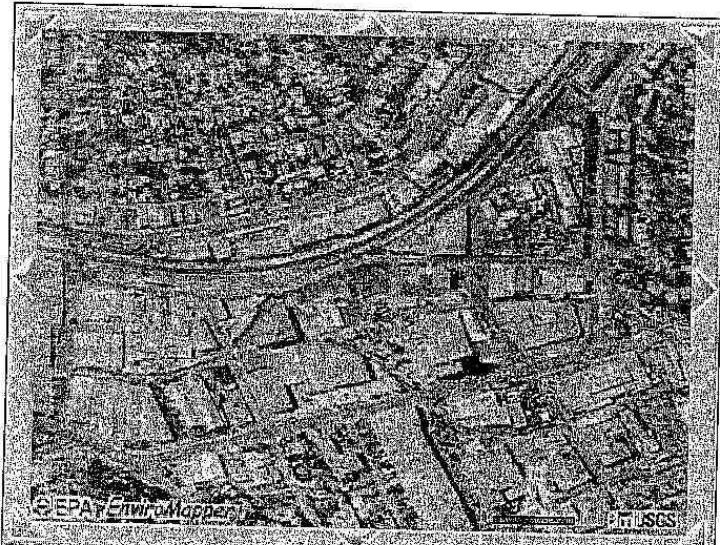
U.S. Environmental Protection Agency Toxics Release Inventory (TRI) Program

[Recent Additions](#) | [Contact Us](#) | [Print Version](#) Search:

GO

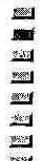
[EPA Home](#) > [Toxics Release Inventory Program](#) > [Reporting Materials](#) > [Facility Siting Tool](#)

- [Browse TRI Topics](#)
- [Get TRI Data](#)
- [Reporting Materials](#)
- [TRI Chemicals](#)
- [TRI PBT Chemicals](#)
- [Laws, Regulations and Executive Order](#)
- [Guidance Documents](#)
- [State TRI Programs](#)
- [International TRI](#)



[Find Another Location](#)

Zoom In



Zoom Out



Latitude:

34°03'43"

Longitude:

-118°10'56"

LEGEND

- Toxic releases
- Cities
- Railroads
- Streets
- Major roads
- Local streets
- Water bodies
- Streams
- States
- Counties

Your goal: Get the center of your facility's production area centered in the map/photo window at the maximum zoom level. This will allow you to get the most precise coordinates.

Zoom-in/out: Select Zoom In, Zoom Out Bar on the right side.

Recenter: When you are not at the maximum zoom-in level, click on the map/photo to center.

Move map/photo: Click the arrow controls around the map/photo window to move in the direction you choose (North, South, East or West or NW, NE, SW, SE).

Mark location: Once you are at the maximum zoom-in level, click the center of the facility production area on the photo. This will mark the coordinates in the Latitude and Longitude fields.

[MAD Code Information](#)

[EPA Home](#) | [Privacy and Security Notice](#) | [Contact Us](#)

Last updated on Tuesday, March 02, 2004

URL: <http://epamap20.epa.gov/tri/emtri.asp>

APPENDIX C: OFFSITE CONSEQUENCE ANALYSIS

