Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8) GreenScreen® for Safer Chemicals (GreenScreen®) Assessment

Prepared for:

State of Washington Department of Ecology

Prepared by:

ToxServices LLC

October 17, 2014



TABLE OF CONTENTS

GreenScreen® Executive Summary for Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8))i
Chemical Name	1
GreenScreen® Summary Rating for Light Aliphatic Petroleum Naphtha	1
Transformation Products and Ratings	
Introduction	2
PhysicoChemical Properties of Light Aliphatic Petroleum Naphtha	3
Group I Human Health Effects (Group I Human)	4
Carcinogenicity (C) Score	4
Mutagenicity/Genotoxicity (M) Score	4
Reproductive Toxicity (R) Score	5
Developmental Toxicity incl. Developmental Neurotoxicity (D) Score	5
Endocrine Activity (E) Score	5
Group II and II* Human Health Effects (Group II and II* Human)	5
Acute Mammalian Toxicity (AT) Group II Score	5
Systemic Toxicity/Organ Effects incl. Immunotoxicity (ST)	5
Group II Score (single dose)	5
Group II* Score (repeated dose)	5
Neurotoxicity (N)	5
Group II Score (single dose)	5
Group II* Score (repeated dose)	6
Skin Sensitization (SnS) Group II* Score	6
Respiratory Sensitization (SnR) Group II* Score	6
Skin Irritation/Corrosivity (IrS) Group II Score	6
Eye Irritation/Corrosivity (IrE) Group II Score	6
Ecotoxicity (Ecotox)	6
Acute Aquatic Toxicity (AA) Score	6
Chronic Aquatic Toxicity (CA) Score	7
Environmental Fate (Fate)	8
Persistence (P) Score	8
Bioaccumulation (B) Score	9
Physical Hazards (Physical)	9
Reactivity (Rx) Score	9
Flammability (F) Score	9
References	10
APPENDIX A: Hazard Benchmark Acronyms	12
APPENDIX B: Results of Automated GreenScreen® Score Calculation for Light Aliphatic Petr Naphtha (CAS# 64742-89-8)	
APPENDIX C: Pharos Output for Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)	

Template Copyright 2014 © Clean Production Action Content Copyright 2014 © ToxServices

APPENDIX D: EPISuite Modeling Results for Light Aliphatic Petroleum Naphtha (CAS #	
8) Sources to Check for GreenScreen® Hazard Assessment	
Licensed GreenScreen® Profilers	19
TABLE OF FIGURES	
Figure 1: GreenScreen® Hazard Ratings for Light Aliphatic Petroleum Naphtha	2
TABLE OF TABLES	
Table 1: Physical and Chemical Properties of Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)	3

GreenScreen® Executive Summary for Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)

Light aliphatic petroleum naphtha is a chemical that functions as a solvent.

Light aliphatic petroleum naphtha was assigned a GreenScreen® Benchmark Score of LT-1 which may be considered equivalent to a Benchmark 1 ("Avoid-Chemical of High Concern") chemical using the full GreenScreen® method as it has High Group I Human Toxicity (carcinogenicity (C) and mutagenicity (M)). This corresponds to a GreenScreen® benchmark classification 1 e in CPA 2011.

Under the scope of this project, ToxServices screened all paint components against Clean Production Action's GreenScreen® List Translator (LT). Those identified as List Translator Benchmark 1 chemicals ("LT-1") do not undergo a full GreenScreen® evaluation to save time and resources. Per the scope of work, only those hazard scores driven by authoritative listings in the List Translator search were to be assigned. Upon inspection of the dataset, ToxServices expanded the assessments for all LT-1 chemicals in order to evaluate aquatic toxicity and environmental fate, as these endpoints are highly relevant to the alternatives assessment of nonbiocide boat paints. The expanded environmental fate and toxicity literature search or modeling for light aliphatic petroleum naphtha did not result in the capture of any additional Benchmark 1 combinations.

GreenScreen® Benchmark Score for Relevant Route of Exposure:

As a standard approach for GreenScreen[®] evaluations, all exposure routes (oral, dermal, and inhalation) were evaluated together, so the GreenScreen[®] Benchmark Score of 1 ("Avoid-Chemical of High Concern") is applicable for all routes of exposure.

GreenScreen® Hazard Ratings for Light Aliphatic Petroleum Naphtha

	Grou	ıp I Hı	ıman				Gro	up II a	Eco	tox	Fa	Fate Physical							
С	M	R	D	E	AT		ST	N		SnS*	SnR*	IrS	IrE	AA	CA	P	В	Rx	F
						single	repeated*	single	repeated*										
Н	Н	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	Н	M	vL	Н	NA	NA

Note: Hazard levels (Very High (vH), High (H), Moderate (M), Low (L), Very Low (vL)) in *italics* reflect estimated values, authoritative B lists, screening lists, weak analogues, and lower confidence. Hazard levels in **BOLD** font are used with good quality data, authoritative A lists, or strong analogues. Group II Human Health endpoints differ from Group II* Human Health endpoints in that they have four hazard scores (i.e., vH, H, M, and L) instead of three (i.e., H, M, and L), and are based on single exposures instead of repeated exposures. Please see Appendix A for a glossary of hazard acronyms. NA: Not assessed.

GreenScreen® Assessment for Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)

Method Version: GreenScreen® Version 1.21

Assessment Type²: Certified

Chemical Name: Light aliphatic petroleum naphtha

CAS Number: 64742-89-8

GreenScreen® Assessment Prepared By: Quality Control Performed By:

Name: Nikki Maples-Reynolds, M.S.

Name: Jennifer Rutkiewicz, Ph.D.

Title: Toxicologist Title: Toxicologist

Organization: ToxServices LLC Organization: ToxServices LLC

Date: October 3, 2014 Date: October 17, 2014

Assessor Type: Licensed GreenScreen® Profiler

Confirm application of the *de minimus* rule³: N/A

Chemical Structure(s):

No structure identified for this complex mixture of hydrocarbons.

Also called: EINECS 265-192-2; Light aliphatic solvent naphtha (petroleum); Solvent naphtha (petroleum), light aliph.; Solvent naphtha, petroleum, light aliph.; Solvent naphtha (petroleum), light aliphatic (ChemIDplus 2014)

Chemical Structure(s) of Chemical Surrogates Used in the GreenScreen®:

Light aliphatic petroleum naphtha is a member of OECD's C7-C9 Aliphatic Hydrocarbon Solvents Category (OECD 2010) and ECHA's Low Boiling Point Naphthas (Gasolines) Category (ECHA 2014). Light aliphatic petroleum naphtha has varying compositions of C7 to C9 paraffins with varying compositions of normal paraffins, isoparaffins, and/or cycloparaffins. Category members are compositionally similar refined petroleum derivatives that are expected to have similar toxicological properties based on structural similarity. Data for members of these categories were used in the environmental toxicity and fate assessment.

Identify Applications/Functional Uses:

1. Solvent

<u>GreenScreen® Summary Rating for Light Aliphatic Petroleum Naphtha</u>⁴: Light aliphatic petroleum naphtha was assigned a GreenScreen® Benchmark Score of LT-1 which may be considered equivalent to

¹ Use GreenScreen® Assessment Procedure (Guidance) V1.2

² GreenScreen® reports are either "UNACCREDITED" (by unaccredited person), "AUTHORIZED" (by Authorized GreenScreen® Practitioner), "CERTIFIED" (by Licensed GreenScreen® Profiler or equivalent) or "CERTIFIED WITH VERIFICATION" (Certified or Authorized assessment that has passed GreenScreen® Verification Program)

³ Every chemical in a material or formulation should be assessed if it is:

^{1.} intentionally added and/or

^{2.} present at greater than or equal to 100 ppm

⁴ For inorganic chemicals with low human and ecotoxicity across all hazard endpoints and low bioaccumulation potential, persistence alone will not be deemed problematic. Inorganic chemicals that are only persistent will be evaluated under the criteria for Benchmark 4.

a Benchmark 1 ("Avoid-Chemical of High Concern") chemical using the full GreenScreen® method as it has High Group I Human Toxicity (carcinogenicity (C) and mutagenicity (M)). This corresponds to a GreenScreen® benchmark classification 1 e in CPA 2011, 2012a. Under the scope of this project, ToxServices screened all paint components against Clean Production Action's GreenScreen® List Translator (LT). Those identified as List Translator Benchmark 1 chemicals ("LT-1") do not undergo a full GreenScreen® evaluation to save time and resources. Per the scope of work, only those hazard scores driven by authoritative listings in the List Translator search were to be assigned. Upon inspection of the dataset, ToxServices expanded the assessments for all LT-1 chemicals in order to evaluate aquatic toxicity and environmental fate, as these endpoints are highly relevant to the alternatives assessment of nonbiocide boat paints. The expanded environmental fate and toxicity literature search or modeling for light aliphatic petroleum naphtha did not result in the capture of any additional Benchmark 1 combinations.

Figure 1: GreenScreen® Hazard Ratings for Light Aliphatic Petroleum Naphtha

Group I Human							•	Gro	up II a	nd II* Hu		Ecotox Fate					Physical			
	C	M	R	D	E	AT		ST	N		SnS*	SnR*	IrS	IrE	AA	CA	P	В	Rx	F
I							single	repeated*	single	repeated*										
	Н	Н	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	Н	M	vL	Н	NA	NA

Note: Hazard levels (Very High (vH), High (H), Moderate (M), Low (L), Very Low (vL)) in *italics* reflect estimated values, authoritative B lists, screening lists, weak analogues, and lower confidence. Hazard levels in **BOLD** font are used with good quality data, authoritative A lists, or strong analogues. Group II Human Health endpoints differ from Group II* Human Health endpoints in that they have four hazard scores (i.e., vH, H, M, and L) instead of three (i.e., H, M, and L), and are based on single exposures instead of repeated exposures. Please see Appendix A for a glossary of hazard acronyms. NA: Not assessed.

Transformation Products and Ratings:

Identify feasible and relevant fate and transformation products (i.e., dissociation products, transformation products, valence states) **and/or moieties of concern**⁵

Transformation products were not assessed, as Light aliphatic petroleum naphtha is an LT-1 chemical and its score will not be impacted by those of transformation products.

Introduction

A complex combination of hydrocarbons obtained from the distillation of crude oil or natural gasoline. It consists predominantly of saturated hydrocarbons having carbon numbers predominantly in the range of C5 through C10 and boiling in the range of approximately 35°C to 160°C (95°F to 320°F) (ECHA 2014).

ToxServices assessed Light aliphatic petroleum naphtha against GreenScreen® Version 1.2 (CPA 2013) following procedures outlined in ToxServices' SOP 1.69 (GreenScreen® Hazard Assessment) (ToxServices 2013).

⁵ A moiety is a discrete chemical entity that is a constituent part or component of a substance. A moiety of concern is often the parent substance itself for organic compounds. For inorganic compounds, the moiety of concern is typically a dissociated component of the substance or a transformation product.

GreenScreen® List Translator Screening Results

The GreenScreen® List Translator identifies specific authoritative or screening lists that should be searched to identify GreenScreen® benchmark 1 chemicals (CPA 2012b). Pharos (Pharos 2014) is an online list-searching tool that is used to screen chemicals against the List Translator electronically. It checks all of the lists in the List Translator with the exception of the U.S. Department of Transportation (U.S. DOT) lists (U.S. DOT 2008a,b) and these should be checked separately in conjunction with running the Pharos query. The output indicates benchmark or possible benchmark scores for each human health and environmental endpoint. The output for Light aliphatic petroleum naphtha can be found in Appendix C and a summary of the results can be found below:

- Carcinogenicity
 - o EC REACH Annex XVII (EU CMR(1)) Category 2 Carcinogen
 - o EU Risk Phrase R45 may cause cancer
 - o GHS Hazard Statement H350 may cause cancer
 - o EC CLP Inventory (EU CMR(2)) Category 1 B Carcinogen
- Gene mutation
 - o EU Risk Phrase R46 may cause heritable genetic damage
 - o GHS Hazard Statement H340 may cause genetic defects
 - o EC CLP Inventory (EU CMR(2)) Category 1 B Mutagen
- Mammalian
 - o GHS Hazard Statement H304 may be fatal if swallowed and enters airway
 - o EU Risk Phrase R65 Harmful: may cause lung damage if swallowed
- PBT
 - o Canada DSL substances that are persistent
- Flammability
 - o Canada CSST WHMIS Class B2 Flammable liquid
- Restricted List
 - ChemSec Substitution List Classified CMR (Carcinogen, Mutagen &/or Reproductive Toxicant)
 - o German FEA (VwVwS) Class 2 Hazard to water
 - o Environment Canada DSL inherently toxic to humans
 - o Environment Canada DSL inherently toxic to environment

PhysicoChemical Properties of Light Aliphatic Petroleum Naphtha

Light aliphatic petroleum naphtha is a liquid petroleum mixture under standard temperature and pressure. It has a vapor pressure of 4-240 hPa indicating that it will likely exist in the vapor and liquid phases. It has a log K_{ow} estimated to be >3, indicating that it is more soluble in octanol than in water and that it has the potential to bioaccumulate in aquatic biota.

Table 1: Physical and Chemical Properties of Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)												
Property Value Reference												
Molecular formula	Mixture - variable											
SMILES Notation	Mixture - variable											
Molecular weight	Mixture - variable											

Table 1: Physical and	Table 1: Physical and Chemical Properties of Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)													
Property	Value	Reference												
Physical state	Liquid ⁶	ECHA 2014												
Appearance	Colorless low viscosity, mobile liquid ⁷	ECHA 2014												
Melting point	<-60°C ⁸	ECHA 2014												
Vapor pressure	4-240 KPa at 37.8°C 9	ECHA 2014												
Water solubility	Not identified													
Dissociation constant	Not identified													
Density/specific gravity	0.62 - 0.88 at 15°C ¹⁰	ECHA 2014												
Partition coefficient	$Log Kow > 3^{11}$	ECHA 2014												

Hazard Classification Summary Section:

Group I Human Health Effects (Group I Human)

Carcinogenicity (C) Score (H, M, or L): H

Light aliphatic petroleum naphtha was assigned a score of High for carcinogenicity based on presence on authoritative lists. GreenScreen[®] criteria classify chemicals as a High hazard for carcinogenicity when the chemical is listed on the EU CMR (1 or 2) lists or is associated with R45 or H350 (CPA 2012a).

- Authoritative and Screening Lists
 - o Authoritative: EC REACH Annex XVII (EU CMR(1)) Category 2 Carcinogen
 - o Authoritative: EU Risk Phrase R45 may cause cancer
 - o Authoritative: GHS Hazard Statement H350 may cause cancer
 - o Authoritative: EC CLP Inventory (EU CMR(2)) Category 1 B Carcinogen
 - o Screening: Not present on any screening lists

Mutagenicity/Genotoxicity (M) Score (H, M, or L): H

Light aliphatic petroleum naphtha was assigned a score of High for mutagenicity/genotoxicity based on presence on authoritative lists. GreenScreen® criteria classify chemicals as a High hazard for mutagenicity/genotoxicity when the chemical is listed on the EU CMR (1 or 2) lists or is associated with R46 or H340 (CPA 2012a).

- o Authoritative: EU Risk Phrase R46 may cause heritable genetic damage
- o Authoritative: GHS Hazard Statement H340 may cause genetic defects

⁶ Read across from Gasoline - Representative substance for the category of low Boiling Point Naphthas

⁷ Read across from Gasoline - Representative substance for the category of low Boiling Point Naphthas

 $^{^8}$ Gasoline (motor and aviation gasoline) has a melting point of < -60°C (read across from Gasoline - Representative substance for the category of low Boiling Point Naphthas)

⁹ The vapor pressures of Low Boiling Point Naphthas (Gasolines) are specified in the range from 4-240 KPa at 37.8°C as measured by EN 13016-1 (read across from Gasoline - Representative substance for the category of low Boiling Point Naphthas)

¹⁰ The density of low boiling point (gasoline) naphthas as measured by EN ISO 12185 lies in the 0.62 - 0.88 range at 15°C (read across from Gasoline - Representative substance for the category of low Boiling Point Naphthas)

 $^{^{11}}$ Most of the hydrocarbon blocks comprising gasoline have a Log $K_{\rm ow} > 3$, indicating these constituents have a potential to bioaccumulate. However, biotransformation is expected to play an important mitigating role in limiting actual bioaccumulation, particularly in higher organisms that possess extensive enzymatic capabilities to metabolize hydrocarbons. However, this endpoint has been calculated for representative hydrocarbon structures using the BCFWIN v2.16 model within EPISuite 3.12 as input to the hydrocarbon block method incorporated into the PETRORISK model (ECHA 2014).

- o Authoritative: EC CLP Inventory (EU CMR(2)) Category 1 B Mutagen
- o Screening: Not present on any screening lists

Reproductive Toxicity (R) Score (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Developmental Toxicity incl. Developmental Neurotoxicity (D) Score (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists

Screening: Not present on any screening lists

Endocrine Activity (E) Score (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Group II and II* Human Health Effects (Group II and II* Human)

Note: Group II and Group II* endpoints are distinguished in the v 1.2 Benchmark system. For Systemic Toxicity and Neurotoxicity, Group II and II* are considered sub-endpoints and test data for single or repeated exposures may be used. If data exist for single OR repeated exposures, then the endpoint is not considered a data gap. If data are available for both single and repeated exposures, then the more conservative value is used.

Acute Mammalian Toxicity (AT) Group II Score (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Systemic Toxicity/Organ Effects incl. Immunotoxicity (ST) Group II Score (single dose) (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Group II* Score (repeated dose) (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Neurotoxicity (N)

Group II Score (single dose) (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Group II* Score (repeated dose) (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Skin Sensitization (SnS) Group II* Score (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Respiratory Sensitization (SnR) Group II* Score (H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Skin Irritation/Corrosivity (IrS) Group II Score (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Eye Irritation/Corrosivity (IrE) Group II Score (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Ecotoxicity (Ecotox)

Acute Aquatic Toxicity (AA) Score (vH, H, M, or L): H

Light aliphatic petroleum naphtha was assigned a score of High for acute aquatic toxicity based on aquatic toxicity values ranging from 3.1 to 64 mg/L. GreenScreen® criteria classify chemicals as a High hazard for acute aquatic toxicity when the most conservative acute aquatic toxicity values range between 1 and 10 mg/L (CPA 2012a).

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Environment Canada DSL inherently toxic to environment
- OECD 2014
 - \circ Classification as inherently toxic on the DSL is based on modeled data. The pivotal value used for the conclusion was 0.05 mg/L.
- ECHA 2014

Note: Data presented are for members of the Low Boiling Point Naphthas (Gasolines) that were identified as read across chemicals in the REACH dossier.

- o An LL₅₀ of 8.2 mg/L was determined in the fish (*Pimephales promelas*, 96-hr).
- o An LL₅₀ of 10 mg/L was determined in the fish (*Pimephales promelas*, 96-hr).
- An LL₅₀ of 11 mg/L was determined in the fish (*Pimephales promelas*, 96-hr).
- o An EL₅₀ of 4.5 mg/L was determined in the aquatic invertebrate (*Daphnia magna*, 48-hr).
- o An EL₅₀ of 12 mg/L was determined in the aquatic invertebrate (*Daphnia magna*, 48-hr).
- o An EL₅₀ of 8.4 mg/L was determined in the aquatic invertebrate (*Daphnia magna*, 48-hr).
- o An EL₅₀ of 3.1 mg/L was determined in algae (*Selenastrum capricornutum*, 72-hr).

- o An EL₅₀ of 3.7 mg/L was determined in algae (*Selenastrum capricornutum*, 96-hr).
- o An EC₅₀ of 64 mg/L was determined in algae (*Selenastrum capricornutum*, 96-hr).
- o An EL₅₀ of 6.4 mg/L was determined in algae (*Selenastrum capricornutum*, 96-hr).

• OECD 2010

- Multi-constituent members of OECD's C7-C9 Aliphatic Hydrocarbon Solvents Category have measured LC/EC₅₀ values from 0.3-1.3 mg/L in fish (96-h), 0.7-0.9 mg/L in daphnia (48-h), and 0.4 mg/L in algae (72-h). The corresponding LL/EL50 values were not provided.
- Since classification on the DSL was based on modeled data, this classification on the screening list was disregarded in favor of measured data. While the acute aquatic toxicity values for low boiling point naphthas ranged from 3.1 to 64 mg/L, the weight of evidence suggests that the acute aquatic toxicity values tend to range between 1 and 10 mg/L, indicating that light aliphatic petroleum naphtha is associated with a High hazard for acute aquatic toxicity. Only LC/EC₅₀ values were provided for OECD's multi-constituent members of OECD's C7-C9 Aliphatic Hydrocarbon Solvents Category. Light aliphatic petroleum naphtha is a complex mixture of compounds, with solubility of each varying from negligible to slight (OECD 2010). The use of LL/EL₅₀ values rather than LC/EC₅₀ values is more appropriate for complex substances with a wide range of water solubilities (UN 2013). Therefore the score of High was assigned based on numerous LL/EL₅₀ values of compositionally similar compounds.

Chronic Aquatic Toxicity (CA) Score (vH, H, M, or L): M

Light aliphatic petroleum naphtha was assigned a score of Moderate for chronic aquatic toxicity based on the weight of evidence suggesting that chronic aquatic toxicity values for light aliphatic petroleum naphtha are generally between 1 and 10 mg/L. GreenScreen® criteria classify chemicals as a Moderate hazard for chronic aquatic toxicity when chronic aquatic toxicity values range between 1 and 10 mg/L (CPA 2012a).

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Environment Canada DSL inherently toxic to environment
- OECD 2014
 - \circ Classification as inherently toxic on the DSL is based on modeled data. The pivotal value used for the conclusion was 0.05 mg/L.
- ECHA 2014

Note: Data presented are for members of the Low Boiling Point Naphthas (Gasolines) that were identified as read across chemicals in the REACH dossier.

- A NOELR of 2.6 mg/L was determined in the fish for mortality (*Pimephales promelas*, 14-d)
- o An LL₅₀ of 5.2 mg/L was determined in the fish (*Pimpehales promelas*, 14-d).
- o A NOELR of 2.6 mg/L was determined in the aquatic invertebrate for reproduction (*Dapnia magna*, 21-d).
- o An EL₅₀ of 10 mg/L was determined in the aquatic invertebrate for reproduction (*Daphnia magna*, 21-d).
- o A NOELR of 16 mg/L was determined in the aquatic invertebrate for mortality (*Daphnia magna*, 21-d).
- \circ An EL₅₀ of >40 mg/L was determined in the aquatic invertebrate (*Daphnia magna*, 21-d).
- o A NOELR of 6.3 mg/L was determined in the aquatic invertebrate for reproduction (*Daphnia magna*, 21-d).
- o A NOELR of 16 mg/L was determined in the aquatic invertebrate for mortality (Daphnia

- magna, 21-d).
- o An EL₅₀ of 13 mg/L was determined in the aquatic invertebrate for reproduction (*Daphnia magna*, 21-d).
- An EL₅₀ of 27 mg/L was determined in the aquatic invertebrate for mortality (*Daphnia magna*, 21-d).
- o A NOELR of 0.5 mg/L was determined in algae for growth rate (*Selenastrum capricornutum*, 72-hr).
- A NOELR of 51 mg/L was determined in algae for cell density (Selenastrum capricornutum, 96-hr).
- o A NOELR of 1.9 mg/L was determined in algae for cell density (*Selenastrum capricornutum*, 96-hr).
- A NOELR of 10 mg/L was determined in algae for growth rate (Selenastrum capricornutum, 72-hr).

• OECD 2010

- A multi-constituent member of OECD's C7-C9 Aliphatic Hydrocarbon Solvents Category (not specified) has a 21-day NOELR of 1.0 mg/L (NOEC of 0.17 mg/L) in daphnia based on reproduction.
- Since classification on the DSL was based on modeled data, this classification on the screening list was disregarded in favor of measured data. While the chronic aquatic toxicity values for low boiling point naphthas ranged from 0.5 to 51 mg/L, the weight of evidence suggests that the chronic aquatic toxicity values tend to range between 1 and 10 mg/L, indicating that light aliphatic petroleum naphtha is associated with a moderate hazard for chronic aquatic toxicity.

Environmental Fate (Fate)

Persistence (P) Score (vH, H, M, L, or vL): vL

Light aliphatic petroleum naphtha was assigned a score of Very Low for persistence as the chemical was classified as readily biodegradable. GreenScreen® criteria classify chemicals as a Very Low hazard for persistence when they are determined to be readily biodegradable and the dominant compartment is water (CPA 2012a).

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Environmental Canada DSL substances that are persistent

• ECHA 2014

- An OECD Guideline 301F ready biodegradability test was performed to assess the biodegradability of light aliphatic petroleum naphtha. The test substance achieved 77% biodegradation in 28 days. Additionally, the authors stated that the test substance met the criteria for ready biodegradability.
- Light aliphatic petroleum naphtha had a measured biodegradation rate of 94% in 25 days.
 No details were provided on the ready biodegradability of the substance, and the authors categorized light aliphatic petroleum naphtha as "inherently biodegradable".
- Light aliphatic petroleum naphtha achieved 74% biodegradation in 28 days. The authors categorized the test substance as "inherently biodegradable". No additional details were provided.

• OECD 2010

 Multi-constituent members of OECD's C7-C9 Aliphatic Hydrocarbon Solvents Category are expected to be readily biodegradable. OECD reports that multi-constituent members demonstrate > 60% degradation after 28 days but do not meet the 10-day window in tests

according to OECD Guideline 301F

- U.S. EPA 2012
 - The BIOWIN modeling Ready Biodegradable Predictor indicates that light aliphatic petroleum naphtha is expected to be readily biodegradable. Fugacity modeling predicts 63.4% will partition to water with a half-life of 9 days, 34.9% will partition to air with a half-life of 1.5 days, 1.19% will partition to soil with a half-life of 17 days, and <1% will partition to sediment with a half-life of 78 days (Appendix D).
- In general, measured biodegradation data indicate that light aliphatic petroleum naphtha is expected to be readily biodegradable. This is supported by two studies that, although not ready biodegradability studies, illustrate that the chemical will biodegrade quickly and OECD's report that multi-constituent category members are expected to be readily biodegradable. Further, biodegradation modeling supports that this chemical will be readily biodegradable. Although the chemical is listed on the DSL as a persistent chemical, no details are available regarding this classification (OECD 2014) and the weight of evidence indicates that this chemical will not be persistent in its dominant compartment (water).

Bioaccumulation (B) Score (vH, H, M, L, or vL): H

Light aliphatic petroleum naphtha was assigned a score of High for bioaccumulation based on an estimated partition coefficient of 4.66. GreenScreen® criteria classify chemicals as a High hazard for bioaccumulation when the partition coefficient is between 4.5 and 5 and BCF values are between 1,000 and 5,000 (CPA 2012a). Confidence was reduced due to lack of measured data for this complex mixture.

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists
- OECD 2010
 - C7-C9 Aliphatic Hydrocarbon Solvents Category have low to moderate potential for bioaccumulation based on an experimental BCF of 199 in mussels for one category member (n-octane) and modeled BCF values of 105-1,216 for constituents.
- U.S. EPA 2012
 - o BCFBAF predicts a BCF of 682 based on a log K_{ow} of 4.66 (Appendix D).
- Modeling predicts a BAF of 682 for the light aliphatic petroleum naphtha mixture, but the structure
 modeled may not represent the chemical components that have the highest bioaccumulation
 potential. Therefore a conservative score of High was assigned based on the estimated log K_{ow} of
 4.66 and the highest estimated BCF of 1,216 for constituents.

Physical Hazards (Physical)

Reactivity (Rx) Score (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Not present on any screening lists

Flammability (F) Score (vH, H, M, or L): Not Assessed

- Authoritative and Screening Lists
 - o Authoritative: Not present on any authoritative lists
 - o Screening: Canada CSST WHMIS Class B2 Flammable liquid

References

ChemIDplus. 2014. Entry for light aliphatic petroleum naphtha (CAS# 64742-89-8). United States National Library of Medicine. Available at: http://chem.sis.nlm.nih.gov/chemidplus/chemidheavy.jsp

Clean Production Action (CPA). 2011. The GreenScreen® for Safer Chemicals Version 1.2 Benchmarks. Dated October 2011. Available at: http://www.greenscreenchemicals.org/

Clean Production Action (CPA). 2012a. The GreenScreen® for Safer Chemicals Version 1.2 Criteria. Dated: November 2012. Available at: http://www.greenscreenchemicals.org/

Clean Production Action (CPA). 2012b. List Translator. Dated February 2012. Available at: http://www.greenscreenchemicals.org/

Clean Production Action (CPA). 2013. The GreenScreen® for Safer Chemicals Chemical Hazard Assessment Procedure. Version 1.2 Guidance. Dated August 31, 2013. Available at: http://www.greenscreenchemicals.org/

European Chemicals Agency (ECHA). 2014. REACH Dossier for light aliphatic petroleum naphtha (CAS# 64742-89-8). Available at: http://apps.echa.europa.eu/registered/data/dossiers/DISS-9ffa5be0-7fdf-37e0-e044-00144f67d031_DISS-9ffa5be0-7fdf-37e0-e044-00144f67d031.html

Grandjean, P. and P.J. Landrigan. 2006. Developmental neurotoxicity of industrial chemicals. Lancet 368: 2167-2178.

Grandjean, P. and P.J. Landrigan. 2014. Neurobehavioral effects of developmental toxicity. The Lancet 13: 330-338.

Organisation for Economic Co-operation and Development (OECD). 2010. SIDS Initial Assessment Profile for C7-C9 Aliphatic Hydrocarbon Solvents Category. SIAM 30, 20-22 April 2010. Available at: http://webnet.oecd.org/Hpv/UI/handler.axd?id=afd8ccb9-af39-43ca-b49c-5034972e75dc.

Organisation for Economic Co-operation and Development (OECD). 2014. Categorization Results from the Canadian Domestic Substance List. Solvent naphtha (petroleum), light aliph. (CAS# 64742-89-8). Available at: http://webnet.oecd.org/CCRWEB/ChemicalDetails.aspx?ChemicalID=23481fe5-5335-43d6-97c0-4259ac05b3b2.

Pharos. 2014. Pharos Chemical and Material Library Entry for light aliphatic petroleum naphtha (CAS# 64742-89-8). Available at: http://www.pharosproject.net/material/

ToxServices. 2013. SOP 1.69: GreenScreen® Hazard Assessments. Dated: August 17, 2013.

United Nations. Globally Harmonized System of Classification and Labelling of Chemicals (GHS) Fifth Revised Edition. Annex 9 Guidance on Hazards to the aquatic environment. Available at: http://www.unece.org/fileadmin/DAM/trans/danger/publi/ghs/ghs_rev05/English/13e_annex9.pdf.

Template Copyright 2014 © Clean Production Action Content Copyright 2014 © ToxServices

United States Department of Transportation (U.S. DOT). 2008a. Chemicals Listed with Classification. 49 CFR § 172.101. Available at: http://www.gpo.gov/fdsys/pkg/CFR-2008-title49-vol2/pdf/CFR-2008-title49-vol2-sec172-101.pdf

United States Department of Transportation (U.S. DOT). 2008b. Classification Criteria. 49 CFR § 173. Available at: http://www.ecfr.gov/cgi-bin/text-idx?c=ecfr&tpl=/ecfrbrowse/Title49/49cfr173_main_02.tpl

United States Environmental Protection Agency (U.S. EPA). 2012. Estimation Programs Interface (EPI) SuiteTM Web, v4.11, Washington, DC, USA. Available at: http://www.epa.gov/opptintr/exposure/pubs/episuitedl.htm.

APPENDIX A: Hazard Benchmark Acronyms (in alphabetical order)

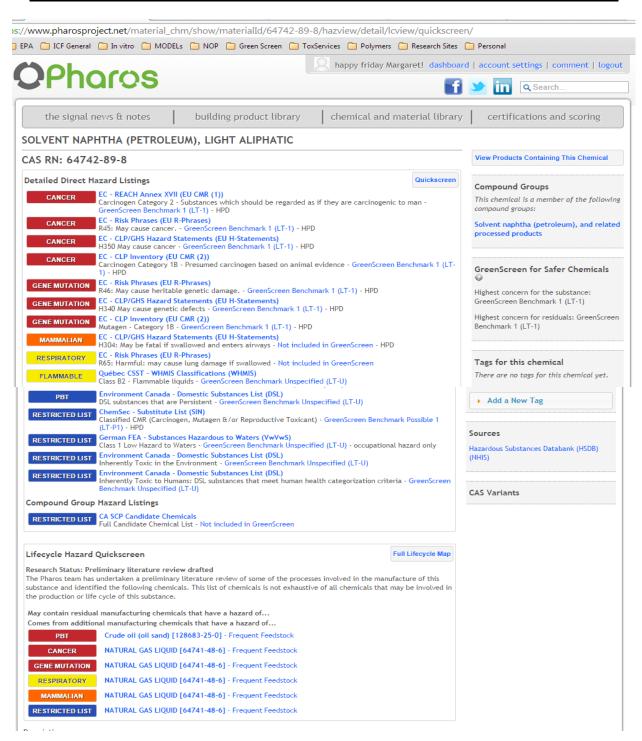
(AA)	Acute Aquatic Toxicity
(AT)	Acute Mammalian Toxicity
(B)	Bioaccumulation
(C)	Carcinogenicity
(CA)	Chronic Aquatic Toxicity
(D)	Developmental Toxicity
(E)	Endocrine Activity
(F)	Flammability
(IrE)	Eye Irritation/Corrosivity
(IrS)	Skin Irritation/Corrosivity
(M)	Mutagenicity and Genotoxicity
(N)	Neurotoxicity
(P)	Persistence
(R)	Reproductive Toxicity
(Rx)	Reactivity
(SnS)	Sensitization- Skin
(SnR)	Sensitization- Respiratory

(ST) Systemic/Organ Toxicity

APPENDIX B: Results of Automated GreenScreen® Score Calculation for Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)

TOXSERVICES									(FreenSo	reen®	Score I	nspecto	r							
TOXICOLOGY RISK ASSE	SSMENT CONSULTING	Table 1: I																	-		
AN SCA				oup I Hun	nan													Fa	ite	sical	
FOR STAFER CHEW			Carcinogenicity Mutagenicity/Genotoxicity Reproductive Toxicity Developmental Toxicity			Endocrine Activity	Acute Toxicity Systemic Toxicity		Neurotoxicity		Skin Sensitization*	Respiratory Sensitization*	Skin Irritation	Eye Irritation	Acute Aquatic Toxicity	Chronic Aquatic Toxicity	Persistence	Bioaccumulation	Reactivity	Flammability	
ical Details								S	R*	S	R*	*	*								
Chemical Name	CAS#	C	М	R	D	E	AT	STs	STr	Ns	Nr	SNS*	SNR*	IrS	IrE	AA	CA	P	В	Rx	F
light aliphatic petroleum naphtha	64742-89-8	Н	Н	DG	DG	DG	DG	DG	DG	DG	DG	DG	DG	DG	DG	Н	М	vL	Н	DG	DG
	1	Table 2. I	Tonoud Con	To	hla							Table 4		1			Table 6		1		
				a	b	c	d	e	f	g			al Name	GreenS	creen®			al Name	Final GreenScreen® Benchmark Score		
				No STOP	No	No	No	Yes				-		1	1		-		1	1	
		3	,	STOP								Note: Chemi	cal has not un				After Data g	ap Assessment ata gap Assessi		Preliminary	
		4		STOP							İ	assessment. I	ot a r mai Gre	cusciceii sc	oic		GS Benchman	rk Score is 1.			
		Table 5: I	Data Gap A	Assessme	nt Table																
				a	b	с	d	e	f	g	h	i	j	bm4	End Result						
												-			1						
		3	3																		
li	cal Details Chemical Name light alliphatic petrole um	Chemical Name ight aliphatic petroleum 64742-89-8	cal Details Chemical Name light aliphatic petroleum naphtha Table 3: I Bench Table 5: I Datagap	Cal Details Chemical Name ight aliphatic petroleum naphtha CAS# C M Benchmark Table 3: Hazard Su Benchmark	cal Details Chemical Name light aliphatic petroleum naphtha CAS# C M R Table 3: Hazard Summary Ta Benchmark a 1 No 2 STOP 3 STOP 4 STOP Table 5: Data Gap Assessme Datagap Criteria a 1 2 3 3	Group I Human April April	Group I Human A	Group I Human State Composition Casternation Casternation	Group I Human Jiji Jiji	Group I Human Jijj Jijj	Group I Human A	Group I Human Group II and IIe	Cal Details		Group I Human Group II and II* Human A tiply A tiply	Group I Human Group II and II* Human Air Joyco	Group I Human Group II and II* Human Lea Details Group Human Group Haman Group Hand He Human Ecotox Air Divisor	Group I Human Group I and II® Human Group II and II® Human Feotox Fi Ai junt Junt Junt Junt Junt Junt Junt Junt J	Group H Human Group H and H* Human Fectors Fate Group H and H* Human Fectors Fate F	Companies Comp	

APPENDIX C: Pharos Output for Light Aliphatic Petroleum Naphtha (CAS# 64742-89-8)



APPENDIX D: EPISuite Modeling Results for Light Aliphatic Petroleum Naphtha (CAS #64742-89-8)

CAS Number: 64742-89-8 SMILES: CCCCCCC CHEM: Solvent naphtha (petroleum), light aliph. MOL FOR: C7 H16 MOL WT: 100.21 ----- EPI SUMMARY (v4.11) -----**Physical Property Inputs:** Log Kow (octanol-water): -----Boiling Point (deg C): -----Melting Point (deg C): -----Vapor Pressure (mm Hg): -----Water Solubility (mg/L): -----Henry LC (atm-m³/mole): -----Log Octanol-Water Partition Coef (SRC): $Log K_{ow} (K_{ow}WIN v1.68 estimate) = 3.78$ $Log K_{ow}$ (Exper. database match) = 4.66 Exper. Ref: MILLER, M.M. ET AL. (1985) Boiling Pt, Melting Pt, Vapor Pressure Estimations (MPBPVP v1.43): Boiling Pt (deg C): 96.16 (Adapted Stein & Brown method) Melting Pt (deg C): -81.01 (Mean or Weighted MP) VP (mm Hg,25 deg C): 45.9 (Mean VP of Antoine & Grain methods) VP (Pa, 25 deg C): 6.12E+003 (Mean VP of Antoine & Grain methods) MP (exp database): -90.6 deg C BP (exp database): 98.5 deg C VP (exp database): 4.60E+01 mm Hg (6.13E+003 Pa) at 25 deg C Water Solubility Estimate from Log K_{ow} (WS K_{ow} v1.42): Water Solubility at 25 deg C (mg/L): 3.554 log K_{ow} used: 4.66 (expK_{ow} database) no-melting pt equation used Water Sol (Exper. database match) = 3.4 mg/L (25 deg C) Exper. Ref: YALKOWSKY, S.H. & DANNENFELSER, R.M. (1992) Water Sol Estimate from Fragments: Wat Sol (v1.01 est) = 2.6374 mg/LECOSAR Class Program (ECOSAR v1.11): Class(es) found: **Neutral Organics** Henrys Law Constant (25 deg C) [HENRYWIN v3.20]: Bond Method: 2.27E+000 atm-m³/mole (2.30E+005 Pa-m³/mole) Group Method: 2.39E+000 atm-m³/mole (2.42E+005 Pa-m³/mole)

Exper Database: 2.00E+00 atm-m³/mole (2.03E+005 Pa-m³/mole) For Henry LC Comparison Purposes: User-Entered Henry LC: not entered Henrys LC [via VP/WSol estimate using User-Entered or Estimated values]: HLC: 1.703E+000 atm-m³/mole (1.725E+005 Pa-m³/mole) VP: 45.9 mm Hg (source: MPBPVP) WS: 3.55 mg/L (source: WSK_{ow}WIN) Log Octanol-Air Partition Coefficient (25 deg C) [K_{0a}WIN v1.10]: Log K_{ow} used: 4.66 (exp database) Log K_{aw} used: 1.913 (exp database) Log K_{oa} (K_{oa}WIN v1.10 estimate): 2.747 Log K_{0a} (experimental database): 2.950 Probability of Rapid Biodegradation (BIOWIN v4.10): Biowin1 (Linear Model): 0.8083 Biowin2 (Non-Linear Model): 0.9686 **Expert Survey Biodegradation Results:** Biowin3 (Ultimate Survey Model): 3.2761 (days-weeks) Biowin4 (Primary Survey Model): 3.9722 (days) MITI Biodegradation Probability: Biowin5 (MITI Linear Model): 0.6619 Biowin6 (MITI Non-Linear Model): 0.8605 Anaerobic Biodegradation Probability: Biowin7 (Anaerobic Linear Model): 0.4892 Ready Biodegradability Prediction: YES Hydrocarbon Biodegradation (BioHCwin v1.01): LOG BioHC Half-Life (days): 0.7385 BioHC Half-Life (days): 5.4758 Sorption to aerosols (25 Dec C)[AEROWIN v1.00]: Vapor pressure (liquid/subcooled): 6.13E+003 Pa (46 mm Hg) Log K_{oa} (Exp database): 2.950 Kp (particle/gas partition coef. $(m^3/\mu g)$): Mackay model: 4.89E-010 Octanol/air (K_{oa}) model: 2.19E-010 Fraction sorbed to airborne particulates (phi): Junge-Pankow model: 1.77E-008 Mackay model: 3.91E-008 Octanol/air (Koa) model: 1.75E-008 Atmospheric Oxidation (25 deg C) [AopWin v1.92]: Hydroxyl Radicals Reaction: OVERALL OH Rate Constant = 6.8713 E-12 cm³/molecule-sec Half-Life = 1.557 Days (12-hr day; 1.5E6 OH/cm³) Half-Life = 18.679 Hrs.Ozone Reaction:

No Ozone Reaction Estimation

Fraction sorbed to airborne particulates (phi):

2.84E-008 (Junge-Pankow, Mackay avg)

1.75E-008 (K_{oa} method)

Note: the sorbed fraction may be resistant to atmospheric oxidation

Soil Adsorption Coefficient (K_{oc}WIN v2.00):

K_{oc}: 239.7 L/kg (MCI method) Log K_{oc}: 2.380 (MCI method)

K_{oc}: 1.107E+004 L/kg (K_{ow} method)

Log K_{oc}: 4.044 (K_{ow} method)

Aqueous Base/Acid-Catalyzed Hydrolysis (25 deg C) [HYDROWIN v2.00]:

Rate constants can NOT be estimated for this structure!

Bioaccumulation Estimates (BCFBAF v3.01):

Log BCF from regression-based method = 2.742 (BCF = 551.7 L/kg wet-wt)

Log Biotransformation Half-life (HL) = 0.2835 days (HL = 1.921 days)

Log BCF Arnot-Gobas method (upper trophic) = 2.834 (BCF = 681.8)

Log BAF Arnot-Gobas method (upper trophic) = 2.837 (BAF = 687.7)

log K_{ow} used: 4.66 (expkow database)

Volatilization from Water:

Henry LC: 2 atm-m3/mole (Henry experimental database)

Half-Life from Model River: 1.022 hours

Half-Life from Model Lake: 95.09 hours (3.962 days)

Removal in Wastewater Treatment (recommended maximum 95%):

Total removal: 99.91 percent Total biodegradation: 0.12 percent Total sludge adsorption: 38.71 percent

Total to Air: 61.08 percent (using 10000 hr. Bio P,A,S)

Removal In Wastewater Treatment (recommended maximum 95%):

Total removal: 99.97 percent

Total biodegradation: 58.86 percent Total sludge adsorption: 17.68 percent

Total to Air: 23.44 percent

(using Biowin/EPA draft method)

Level III Fugacity Model:

Mass Amount Half-Life Emissions

(kg/hr.) (percent) (hr.) 34.9 35.9 1000 Air Water 63.4 208 1000 Soil 1.19 1000 416 Sediment 0.553 1.87e+003 0

Persistence Time: 77 hr.

Sources to Check for GreenScreen® Hazard Assessment

Note: For a GreenScreen[®] Hazard Assessment, data queries should be initially limited to the following references. If data gaps exist after these references have been checked, additional references may be utilized.

U.S. EPA High Production Volume Information System (HPVIS): http://www.epa.gov/hpvis/index.html

UNEP OECD Screening Information Datasets (SIDS): http://www.chem.unep.ch/irptc/sids/OECDSIDS/sidspub.html

OECD Existing Chemicals Database: http://webnet.oecd.org/hpv/ui/SponsoredChemicals.aspx

European Chemical Substances Information System IUCLID Chemical Data Sheets: http://esis.jrc.ec.europa.eu/index.php?PGM=dat

National Toxicology Program: http://ntp.niehs.nih.gov/

International Agency for the Research on Cancer: http://monographs.iarc.fr/ENG/Classification/index.php

Human and Environmental Risk Assessment (HERA) on ingredients of household cleaning products: http://www.heraproject.com/RiskAssessment.cfm

European Chemicals Agency (ECHA) REACH Dossiers: http://echa.europa.eu/

Licensed GreenScreen® Profilers

Light aliphatic petroleum naphtha GreenScreen® Evaluation Prepared by:

Nikki Maples-Reynolds, MS.

Toxicologist

ToxServices LLC

Light aliphatic petroleum naphtha GreenScreen® Evaluation QC'd by:

Jennifer Rutkiewicz, Ph.D.

Toxicologist

ToxServices LLC