SKC Passive Samplers

New Designs for New Applications







VOC Chek 575 and ULTRA ppb to ppm-Level VOCs

See pages 74-75 and 72-73



U.F.

Passive Samplers for Health Care Ethylene Oxide, Isoflurane, Halothane

See page 83

UME^X 100 Formaldehyde/Other Aldehydes

See page 84

UME^X 200 Nitrogen Dioxide/Sulfur Dioxide

See page 85

UME^X 300 Ammonia

See page 85

Elemental Mercury and Hydrogen Cyanide

See page 86

Diffusive TD Tubes BTEX/Other VOCs

Diffusive Thermal Desorption Tubes Time-averaged Assessment of BTEX and Other VOCs

- Validated sampling rates for benzene, toluene, ethylbenzene, xylenes, and many other VOCs
- Contain pre-conditioned Carbopack[®] X sorbent
- SilcoNert[®] 2000 deactivated 3.5 x 0.25-inch OD stainless steel tubes
 Ideal for protecting sample integrity over long sampling periods
- Use to determine average concentration over longer sampling periods up to 14 days
- No pump or training required
- Reusable to keep costs low
- Unique identification number and bar code
- Weatherproof tube shelter available
- Airflow and cap placement arrow indicators printed on tube
- Sealed with brass Swagelok screw caps with PTFE ferrules for transport
- Each collected compound can be analyzed quantitatively

SKC Diffusive Thermal Desorption (TD) Tubes are designed to provide accurate sampling of VOCs at refinery fencelines and for other ambient air monitoring applications without the use of a sample pump. Sampling times of up to 14 days provide an average concentration that eliminates the effects of daily variables such as hazard concentration and temperature. A weatherproof shelter accessory protects samples.

Description	Cat. No.	Qty.
SKC Diffusive TD Tubes,* 3.5 x 0.25-in OD deactivated stainless steel		
tubes filled with pre-conditioned Carbopack X and supplied with diffusion		
caps and brass Swagelok screw caps with PTFE ferrules	226-520	10
Diffusion Caps	226-525	10
Shelter	226-526	ea

* Tubes must be used within 30 days of conditioning

Diffusive TD Tubes BTEX/Other VOCs

SKC Diffusive Thermal Desorption Tubes for BTEX and Other VOCs



SKC Diffusive TD Tubes are conditioned and quality control tested to assure low background.

More Information McClenny, W.A. et al., "24 h Diffusive Sampling of Toxic VOCs in Air onto Carbopack X ...," http://doi.org/fhgn4s U.S. EPA Method 325B: https://goo.gl/ B4viuV www.skcinc.com



ULTRA for Sub-ppb Level VOCs

ULTRA Passive Samplers

Convenient Alternative to Canisters and Thermal Desorption Tubes

Compare ULTRA to Canisters

Side-by-side studies using ULTRA® Passive Samplers (Anasorb GCB1) and stainless steel canisters demonstrate excellent sampling correlation.

Compound	ULTRA (µg/m³)	Canister (µg/m ³)
Benzene	4.2	4.5
	2.1	2.0
	1.9	1.6
	6.67	6.8
	1.58	1.5

Compound	ULTRA (µg/m³)	Canister (µg/m ³)
Perchloro- ethylene	1.1	1.6
	2.3	2.2
	32.9	30.0
	1.37	2.0
	2.85	2.6

Compound	ULTRA (µg/m³)	Canister (µg/m ³)
o-Xylene	7.55	7.9
	1.16	0.93
	1.96	1.9
	8.3	6.2
	13.3	11.0

Compound	ULTRA (µg/m³)	Canister (µg/m ³)
Toluene	30.0	26.0
	20.3	19.0
	44.0	46.0
	10.8	8.8
	6.1	3.8

Compound	ULTRA (µg/m³)	Canister (µg/m ³)
m,p-Xylene	21.2	19.2
	5.52	5.6
	34.1	36.7
	3.7	2.51
	5.7	5.1

Results comparable to canisters for EPA Method TO-15

- No cleaning or certification costs
- Lower purchase price
- No expensive shipping

Passive alternative for EPA TO-17 — no pump required

Choose from 5 sorbents for indoor and ambient air sampling (including SVOCs)

- Anasorb GCB1
- Tenax TA
- Chromosorb 106
- Carbopack X
- Charcoal (solvent extraction)
- Easy on/off sampling
- Small, lightweight, and easy to transport
- Validated sampling (uptake) rates
 - See www.skcinc.com/catalog/passive-guide.php
- Built-in blank/correction sorbent section available
- Sample integrity
 - Manufactured in an ultra-clean environment
 - Extensive cleaning and QC procedures
 - Sonically welded housing
- Long-term sampling up to 30 days reduces temporal variability (see www.skcinc.com/pdf/1812.pdf)
- Ideal for vapor intrusion sampling
- See EPA-OSWER technical guide at www.epa.gov, search "vapor intrusion"



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<section-header>

Ordering Select from prefilled samplers or separate sampler housing and sorbent vials.	Economy Prefilled Samplers, without built-in blank, pk/5	Prefilled Samplers, with built-in blank, pk/5	Sorbent Vials for User-filled Samplers, pk/2, require empty housing Cat. No. 690-200			
Sorbent/Amount	Cat. No.	Cat. No.	Cat. No.			
Anasorb GCB1,## 370 mg in each compartment or vial	690-101-NB	690-101	690-201			
Chromosorb 106,# 285 mg in each compartment or vial	690-103-NB	690-103	690-203			
Tenax TA,# 253 mg in each compartment or vial	690-104-NB	690-104	690-204			
Charcoal,# 500 mg in each compartment or vial (solvent extraction)	690-105-NB	690-105	690-205			
Carbopack X,# 400 mg in each compartment or vial	690-106-NB	690-106	690-206			
Empty Sampler Housing, for user-filled ULTRA, required 690-200						
# Limited shelf-life	o www.osha.gov and search "ULTF	RA" for additional information on sai	mpling rates for Chromosorb 106.			

Sampling Accessories	Cat. No.
Rate Reducer, 12 holes, lowers sampling rate for extended	
sampling time and higher concentrations	
	690-300
Transfer Funnel, for filling sampler housing with sorbent from vials, for ULTRA only	690-301
Stand for Indoor Sampling	
	690-302
Shelter for Outdoor Sampling	
	690-303
Analysis Accessories	Cat. No.
hermal Desorption Tube, Perkin Elmer, 6.35 x 88.9 mm (0.25 x 3.5 inches), includes screens and end caps	P226530
Analysis Transfer Funnel, facilitates transfer of sorbent from vial to 6.35-mm (0.25-inch) OD thermal desorption tube	590-264
U.S. Patent No. 6.607.581	

See www.skcinc.com/catalog/passive-guide.php for sampling rates for over 50 compounds.

SKC 575 Series Passive Samplers are now...



575 Series

Decrease sampling costs

- More sorbent/capacity + validation to 2x exposure level
 - = One sampler per 8-hour shift
 - = One analysis of a single layer
 - = Savings for you

✓ Increase convenience

- No assembly required for sampling
- New design for easier sorbent transfer during analysis
- Reference guides available for laboratories

✓ Increase reliability

- Sampling rates for over 300 compounds (see pages 76-82)
- Online research reports document performance for legal and compliance issues. Visit www.skcinc.com/reports.asp
- Specialty samplers available for ethylene oxide (EtO), styrene, and methanol
- Validated in OSHA methods

Easy 3-step Sampling!



Remove front cap.





Clip in worker's breathing zone to sample.

Cap and record information.

Passive Samplers

for ppm-Level Organic Vapors



Back of VOC Chek 575 Series Sampler designed for easier analysis

Front of VOC Chek 575 Series Sampler (actual size)

Active/Passive Sampling Cross-reference

Compound	Active Method	Tube Cat. No.	Passive Method	Sampler Cat. No.
Benzene	OSHA 1005	226-01	OSHA 1005	575-002
Butyl acetates	OSHA 1009	226-01	OSHA 1009	575-002
MEK/MIBK	OSHA 16 (MEK)	226-10	OSHA 1004	575-002
Styrene	OSHA 89	226-73	OSHA 1014	575-006
Toluene	OSHA 111	226-81A or 226-01	OSHA 111	575-002
Trichloroethylene/ tetrachloroethylene	OSHA 1001	226-01	OSHA 1001	575-002
Xylenes/ethylbenzene	OSHA 1002	226-01	OSHA 1002	575-002

VOC Chek 575 Series Passive Sampler Ordering

Passive Sampler for:	Sorbent		Cat. No.
Organic vapors	Charcoal Lot 2000, 350 mg	pk/5	575-001 [†]
	-	pk/25	575-001A
Organic vapors	Anasorb 747, 500 mg	pk/5	575-002 [†]
	-	pk/25	575-002A
Ethylene oxide	Anasorb 747 treated with hydrobromic acid, 500 mg	pk/5	575-005
		pk/25	575-005A
Styrene	Anasorb 747 treated with tert-butyl catechol, 500 mg	pk/5	575-006
Methanol	Anasorb 747, 500 mg, includes secondary diffusion barrier	pk/5	575-007

† Larger quantity packages are available. Contact SKC.

SKC Inc. www.skcinc.com



Leader in Passive Sampler Technology

SKC VOC Chek® 575 Series Passive Samplers are identified as a reliable alternative to active sampling for many compounds in OSHA diffusive methods!

Methanol Passive Sampler Cat. No. 575-007 see ordering below left

STOP *No need to double your costs!* VOC Chek Series samplers have: - More sorbent in a

- single layerMore capacity
- Wore capacity
- Validated results to 2x exposure limit
 Lower price than
- double-layer samplers

See the VOC Chek 575 Series Selection Guide pages 76-82

VOC Chek 575 Series Selection Guide On-worker ppm Level Sampling of Organic Vapors

SKC VOC Chek 575 Series Passive Samplers are available with sorbents such as activated carbon and Anasorb 747 that strongly retain the collected compounds and require solvent extraction for laboratory analysis.

Use the following guide to locate target compounds and get an overview of critical sampling parameters along with the SKC catalog number of the recommended SKC VOC Chek 575 Series Passive Sampler. Visit www.skcinc.com and download our free SKC Sampling Guide app to access this information on your Android or Apple device.

Validation Levels

See the Validation Level column in this guide to determine the level of scientific testing. Go to www.skcinc.com and search the number in the Research Report column to access online research reports.

Full – Stringent NIOSH testing protocol has been applied to verify sampler desorption efficiency, sampling rate, capacity, and the effects of relative humidity, temperature, concentration, reverse diffusion, and storage on accuracy.

Bi-level – NIOSH testing protocol has been applied to the most volatile member of a related (homologous) series of chemicals; therefore, related less volatile series members require only partial validation (*described below*) to achieve the same level of sampling accuracy. See Guild et al. (http://doi.org/cjzqd7) or request a copy from SKC.

Partial - NIOSH testing protocol has been applied to verify sampling rate, desorption efficiency, and storage stability.

Calculated – This calculation of a sampling or uptake rate uses the diffusion coefficient of a specific chemical (D) and the cross-sectional area (A) and length (L) of the sampler's diffusion path (based on sampler geometry).

OSHA - Sampler has been validated by OSHA and is referenced in a published OSHA method.

			OSHA	PELs Δ		Sampling Time						
	Validation	Research	TWA	CLG/STEL	Sampling Rate	Min	Max	Analytical	DE %	SKC VC	C Che	k 575
Chemical Hazard	Level	Report	(ppm)	(ppm)	(ml/min)	(min)	(hrs)	Method	§	Cat. No		
Acetic acid	Calculated		10	15 #	19.6			GC-FID	99.2	575-001		
Acetic acid	Calculated		10	15 #	19.6			GC-FID	107.9	575-002		
Acetoin (acetyl methyl carbinol)	Calculated				14.9			GC-FID		575-001	or	575-002
Acetone	Full	1303	1000		20.3 🗸	15	4	GC-FID	90.2	575-002		
Acetone	Full	1303	1000		15.2 ≈	240	8 ≈	GC-FID	90.2	575-002		
Acetonitrile	Calculated		40		22.4			GC-FID	103	575-001		
Acetonitrile	Calculated		40		22.4			GC-FID	108	575-002		
Acetyl methyl carbinol (acetoin)	Calculated				14.9			GC-FID		575-001	or	575-002
Acrylonitrile	Full		2	10 C	20.4	15	8	GC-FID	81	575-002		
Allyl alcohol	Calculated		2	4 #	18.4			GC-FID	64	575-001		
Allyl alcohol	Calculated		2	4 #	18.4			GC-FID	76	575-002		
Allyl chloride	Calculated		1	2 #	17.8			GC-FID	95.1	575-001		
Allyl chloride	Calculated		1	2 #	17.8			GC-FID	101.3	575-002		
n-Amyl acetate	Calculated		100		11.7			GC-FID	93.5	575-001		
n-Amyl acetate	Calculated		100		11.7			GC-FID	96	575-002		
sec-Amyl acetate (2-pentyl acetate)	Calculated		125		11.8			GC-FID		575-001		
n-Amyl alcohol	Calculated				13.9			GC-FID	87.3	575-001		
n-Amyl alcohol	Calculated				13.9			GC-FID	100.6	575-002		
t-Amyl methyl ether (methyl tert-amyl ether)	Bilevel	1355			13.1	30	8	GC-FID	99	575-001		
Aniline	Calculated		5		14.2			GC-FID		575-001		
Benzene	Full	1312	1	5	16	15	8	GC-FID	93.5	575-001		
Benzene	OSHA 1005		1	5	17.1	15	8	GC-FID	93.6	575-002		
Benzotrifluoride (trifluoromethyl benzene; OXSOL 2000)	Bilevel		100		13.3	15	8	GC-FID	106	575-001		
Benzotrifluoride (trifluoromethyl benzene; OXSOL 2000)	Bilevel		100		13.3	15	8	GC-FID	107	575-002		
Benzyl acetate	Calculated		10		11.3			GC-FID	91.2	575-002		
Benzyl chloride	Calculated		1	1 C	12.9			GC-FID	98.7	575-001		
Benzyl chloride	Calculated		1	1 C	12.9			GC-FID	98.9	575-002		
Bromobenzene	Calculated				13.8			GC-FID		575-001		
Bromodichloromethane (dichlorobromomethane)	Calculated				16.1			GC-FID		575-001		
Bromoethane (ethyl bromide)	Calculated		200	250 #	18.5			GC-FID		575-001		
Bromoform	Calculated		0.5		15.2			GC-FID		575-001		
Bromomethane (methyl bromide)	Calculated			20 C	22.1			GC-FID		575-001		
1-Bromopropane (propyl bromide)	Full	1740	0.1		14.4	30	8	GC-FID	100	575-001		
1-Bromopropane (propyl bromide)	Full	1740	0.1		14.7	30	8	GC-FID	107	575-002		
1,3-Butadiene	Calculated		1	5	19.6		-	GC-FID		575-001		
n-Butane	Calculated				18.1			GC-FID		575-001		
n-Butanol (n-butyl alcohol)	Calculated		100	50 #	15.5			GC-FID	94	575-001		

See page 212 for abbreviations.

Value <th< th=""><th></th><th></th><th></th><th>OSHA</th><th>PELs 🛆</th><th></th><th colspan="2">Sampling</th><th></th><th></th><th colspan="2"></th></th<>				OSHA	PELs 🛆		Sampling					
Distance Calaxies Park	Chamical Hozard	Validation	Research	TWA	CLG/STEL	Sampling Rate	Min (min)	Max	Analytical	DE %	SKC VOC Chek 575	
Seared loss-out activity Column of the second activity of the second activ	n-Butanol (n-butyl alcohol)	Calculated	нероп	(ppm) 100	(ppm) 50 #	(m/min)	(min)	(nrs)	GC-FID	<u>s</u> 100	575-002	
Based scalarsBased and and and and and and and and and an	2-Butanol (sec-butyl alcohol)	Calculated		150	150 #	15.6			GC-FID	93	575-001	
Balance instrying Bited Size Join International status Size Size Balance instrying Caluate Ro Size Si	2-Butanol (sec-butyl alcohol)	Calculated		150	150 #	15.6			GC-FID	100	575-002	
Balance Bit Above Control Bit Above Bi	2-Butanone (methyl ethyl ketone, MEK)	Bilevel	1306	200		17.1	15	12	GC-FID	100	575-002	
Base And Lates Calculated No. L2 Co-PPO R2 SP-002 Base And Lates Calculated Calculated SP-00 R2 SP-002 PC-PO R2 SP-002 Base And Lates Calculated Calculated Z20 V R2 SP-002 V R2 <	2-Butanone (methyl ethyl ketone, MEK)	OSHA 1004		200		16.88		8	GC-FID	92.3	575-002	
Baly access OSM 100 158 20.7 5.0 8 67.00 98.2 75.30 see Day locates Crickler 1964 200 203 20 6 67.00 98.2 75.00 see Day locates Crickler 200 12.3 10 8 67.00 98.2 75.00 see Day locates Crickler 200 12.7 15 8 67.40 98.1 75.00 See Day locates Crickler 200 12.7 1 6 67.40 98.1 75.00 See Day locates Crickler 200 12.7 1 6 67.40 98.1 75.00 1 12.00	2-Butoxyethanol (butyl CELLOSOLVE solvent)	Calculated		50		12			GC-FID	89.7	575-002	
Party Journal Party	n-Butyl acetate	OSHA 1009	4004	150	200 #	13.07	15	8	GC-FID	99.2	575-002	
Date data data One Oot Oot Oot Oot <	n-Butyl acetate	Partial	1894	150	200 #	12.3	30	8	GC-FID	90.4	575-002	
See Bulk status Calculator 200 127 15 8 CP-00 89.00 27.9400 Bigl status Calculator Calculator 200 12.7 15 8 CP-00 88.0 75.400 Bigl status Calculator Calculator 200 12.7 15 8 CP-00 88.0 75.400 Bigl status Calculator 200 12.00 15 8 CP-00 86.0 75.400 Status Calculator 100 59.0 15.5 - CP-00 160 75.400 Status Calculator 100 59.0 15.5 - CP-00 100 75.400 Status Calculator 100 59.0 15.5 - CP-00 100 75.400 Status Calculator 100 59.0 55.2 11.3 CP-00 100 75.400 Status Calculator 100 75.00 100 75.400 100	sec-Butyl acetate	Calculated	1034	200	200 #	12.8		0	GC-FID	96.2	575-001	
Description Object voide (expl voide) Constraint Object voide) Constraint Constraint <thconstraint< th=""> <thconstraint< th=""> <thc< td=""><td>sec-Butyl acetate</td><td>Calculated</td><td></td><td>200</td><td></td><td>12.8</td><td></td><td></td><td>GC-FID</td><td>96.6</td><td>575-002</td></thc<></thconstraint<></thconstraint<>	sec-Butyl acetate	Calculated		200		12.8			GC-FID	96.6	575-002	
Bay isotate Cancel of Canc	sec-Butyl acetate	OSHA 1009		200		12.74	15	8	GC-FID	98.9	575-002	
eBuy actation Calculated Obi N4 1009 12.7 Image: Construct on the second of the	t-Butyl acetate	Calculated		200		12.7			GC-FID	95.1	575-001	
Bulk resulta ORHA 100 200 11.0 10 8 6 CPID 88.0 75:002 Bulk result 100	t-Butyl acetate	Calculated		200		12.7			GC-FID	94.8	575-002	
Bayle action Binerel 101 117 30 8 GC+P0 85 75 62 Bayle action Canculated 100 100 15.5 IC GC+P0 85 75 76	t-Butyl acetate	OSHA 1009		200		13.09	15	8	GC-FID	98.9	575-002	
Carbon Carbon Carbon Fig. 2 Control Control <thcontrol< th=""> <thcontrol< th=""> <thcontr< td=""><td>Butyl acrylate</td><td>Bilevel</td><td></td><td>10 ‡</td><td>450</td><td>11.7</td><td>30</td><td>8</td><td>GC-FID</td><td>95</td><td>575-002</td></thcontr<></thcontrol<></thcontrol<>	Butyl acrylate	Bilevel		10 ‡	450	11.7	30	8	GC-FID	95	575-002	
Hardy Action (* Natural) Deckname 100 20 * 132 DCPU 100 55002 Body Body (* Natural) Education 100 <	t-Butyl alcohol	Calculated		100	150	15.8			GC-FID	84	575-002	
Base Burgh action (2) Evaluation) Calculated 150 156 0 62-PD 88 97501 Bayl action (2) Evaluation) Calculated 150 150 156 0 62-PD 88 97501 Bayl action (2) Evaluation Calculated 5 10.4 8 62-PD 975-022 Bayl action (2) Evaluation Calculated 5 10.4 8 62-PD 101 575-012 Bayl action (2) Evaluation Bilevel 150 5 8 62-PD 101 575-012 Bayl action (2) Evaluation Bilevel 10 10.4 15 8 62-PD 103 575-002 Bayl action (2) Evaluation Calculated 11.3 62-PD 103 575-002 Bayl action (2) Evaluation Calculated 2 mpmin 10.8 62-PD 103 575-002 Calculated (2) Evaluation Calculated (2) Empinity Calculated (2) Empinity 10.8 62-PD 135 8 62-PD 135 8 62-PD	n-Butyl alcohol (1-butanol)	Calculated		100	50 #	15.5			GC-FID	94	575-002	
Books (Extund) Calculated 150 150 150 CP-D 100 255/02 Book (EXCOX)/E control (Free dyot monotory letter acted) Calculated 5 10.4 CP-D 755/02 Book (EXCOX)/E control (Free dyot monotory letter acted) Calculated 5 10.4 15 6 CP-D 755/02 Book (GVC) (Free dyot end on the control (Free dyot end on the contro	sec-Butyl alcohol (2-butanol)	Calculated		150	150 #	15.6			GC-FID	93	575-001	
Bayl Bergene encode Calculated Form ST-500 ST-500 Buly div (blue) Clue) Clue Claubide 5 13.1 15.8 6 C-FD 17.500 Buly div (blue) Clue) Clue Claubide 50 55.8 11.6 8 6 C-FD 10 175.002 Buly divel (blue) Clue) Clue Claubide 50 55.8 11.6 8 6 C-FD 10.0 175.002 Buly divel (blue) Clue) Claubide 11.2 15.8 8 6 C-FD 10.0 175.002 Buly divel (blue) Clue) Claubide 11.3 Claubide 11.3 Claubide 11.3 Claubide Claubide 11.3 Claubide Claubide 11.3 Claubide Claubide Claubide Claubide 2 Claubide	sec-Butyl alcohol (2-butanol)	Calculated		150	150 #	15.6			GC-FID	100	575-002	
Biker S 10.4 S CPD 57.002 Biker Biker Biker 135 S	t-Butyl benzene	Calculated				11.3			GC-FID		575-002	
Buy i ethy i e	Butyl CELLOSOLVE acetate (ethylene glycol monobutyl ether acetate)	Calculated		5		10.4			GC-FID		575-002	
Buly goody eller Ge-HD INU S75-001 S75-001 bally namby dent (marthy Houng latent, MTBE) Billweil 10 10.4 15.8 8 GC-HD 7.6 75-001	t-Butyl ethyl ether (ethyl tert-butyl ether)	Bilevel	1356			13.1	15	8	GC-FID	101	575-001	
Bally interfly there (melly it buy deter, MTBE) Full 132 132 135 8 8 6C-FLD 101 575-01 mBuly berzene Calculated 11.23 6C-FLD 103 575-001	n-Butyl glycidyl ether	Calculated		50	5.6‡	11.6			GC-FID	104	575-002	
PetH-Bu/Motivation Direct 10 10.4 10.4 10.8 6 GC+D0 10.0 p/short Bee/Bu/Motrane Calculated 11.3 GC+D0 10.8 GC+D0 GC-D0	t-Butyl methyl ether (methyl t-butyl ether, MTBE)	Full	1352			13.6	8.5	8	GC-FID	97.4	575-001	
Productative Calculated Int.3 Corr Corr aprime-Brynolactore Calculated Int.3 Corr S75-002 aprime-Brynolactore Calculated Z mpim ¹ I0.8 Corr S75-002 Camphor Calculated Z mpim ¹ I0.8 Corr G-PID S75-002 Camphor Calculated Z mpim ¹ I0.8 Corr G-PID S75-002 Cathol acid (phent) Calculated Z mpim ¹ I0.8 G-PID S75-001 S75-002 Cathol intrachivide Calculated 100 12.1 G-PID S75-002 S75-002 Collable acid (phent) Calculated 100 12.1 G-PID S75-002 S75-002 Collable acid (phent) Calculated 100 12.1 G-PID S75-002 S75-002 Collable acid (phent) Earlow 13 15 8 G-PID 195 S75-002 Collow-Chromothylibenzene (panchloroblenzont/liuoride; Bilevel 25 0 11.8 15	p-tert-Butyl toluene	Bilevel		10		10.4	15	8	GC-FID	100	575-001	
Description Description <thdescription< th=""> <thdescription< th=""></thdescription<></thdescription<>	sec-Butylbenzene	Calculated				11.23			GC-FID	103	575-002	
Campion Calculated 2 mg/m1 10.8 Calculated 2 mg/m1 Campion Calculated 2 mg/m1 10.8 GC-PID 94.2 575.001 Carbon and globel and	gamma-Butyrolactone	Calculated				16.6			GC-FID	80.9	575-002	
Campion Calculated 2 mg/m End bit Set 15.6 C 14.5 L SC-PD 13 57:001 or 57:002 Carbon add unlide Calculated 5 15.6 C 14.5 IA	Camphor	Calculated		2 mg/m ³		10.8			GC-FID	94.2	575-001	
Catolic and (pheno) Calculated 5 15.6 C 14.5 C BC-FID 875-001 or 575-002 Catoon disulfide Calculated 20 30 19.54 C CC-FID 98.3 575-001 2-CELLOSOLVE acetate (2-ethoxyethyl acetate) Calculated 100 12.1 2 CC-FID 98.3 575-001 2-CELLOSOLVE acetate (2-ethoxyethyl acetate) Calculated 50 ± 13 15 8 GC-FID 98.4 575-002 C-Choreetherbyl tenzene (monochorotoluene; OXSOL 10) Bilevel 50 ± 11.8 15 8 GC-FID 97.5 75.001 CAISOL 100 Calculated 50 ± 11.8 15 8 GC-FID 97.5 75.001 CAISOL 100 Bilevel 25 ° 11.8 15 8 GC-FID 97.5 75.001 CAISOL 100 Bilevel 75 14.41 15 8 GC-FID 97.5 75.001 Chorobacharene Parial 175<	Camphor	Calculated		2 mg/m ³		10.8			GC-FID	113	575-002	
Carbon disultifie Calculated 20 30 19.54 CC-FID 575-001 2-CEL/DSQLVE actain (2-ethoryethyl actate) Calculated 100 25 14.1 30 8 6C-FID 98.4 575-001 1-Chloro-2-methyl berzare (monochlorotoluene; OXSOL 10) Bilevel 50 ‡ 13 15 8 6C-FID 91.8 575-001 1-Chloro-4-methyl berzare (monochlorotoluene; OXSOL 10) Bilevel 50 ‡ 11.8 15 8 6C-FID 91.8 575-001 1-Chloro-4-methyl berzare (monochlorotoluene; OXSOL 10) Bilevel 25 ° 11.8 15 8 6C-FID 91.8 575-001 1-Chloro-4-methyl berzare (monochlorotoluene; OXSOL 10) Bilevel 25 ° 11.8 15 8 6C-FID 87.5 60.0 10.2 575-001 1-Chloro-4-printely Balevel 200 18.3 6 6C-FID 87.5 60.0 8 6C-FID 57.5 60.0 8 6C-FID 57.5 60.0 8 6C-FID 57.5	Carbolic acid (phenol)	Calculated		5	15.6 C	14.5			GC-FID		575-001 or 575-002	
Carbon transhoride Bilevel 10 25 C 14.1 30 8 GC-FID 98.3 575-001 CELLOSOLVE carbital (2-ethoxyethyl acetate) Calculated 100 12.1 GC-FID 95.4 575-002 1C-bitor-2-methyl berzzene (monochrotoluene; CXSOL 10) Bilevel 50 ‡ 13 15 8 GC-FID 91.8 575-002 1C-bitor-4-fiftilyoromethyl/berzzene (parachlorobenzotrifluoride; CNIon-4-fiftilyoromethyl/berzzene Bilevel 25 ° 11.8 15 8 GC-FID 10 575-002 CNIOn-4-fiftilyoromethyl/berzzene (parachlorobenzotrifluoride; CNIon-4-fiftilyoromethyl/berzzene Partial 183 75 14.41 15 8 GC-FID 157:001 CNIOn-4-fiftilyoromethyl/berzzene Partial 183	Carbon disulfide	Calculated		20	30	19.54			GC-FID		575-001	
C4-ELCSUVE acetate (2-ethoxyettry acetate) Cacuated 100 12.1 GC-FID 95.4 p3-402 1-Chtoro-2-mettry berazene (monochtorotabuere, XSOL 10) Bilevel 50 ± 13 15 8 GC-FID 91 575-002 1-Chtoro-2-mettry berazene (monochtorotabuere, XSOL 10) Bilevel 25 0 11.8 15 8 GC-FID 91 575-002 1-Chtoro-4-tritutoromethylberazene (parachtorobenzotrifluoride; Bilevel 25 0 11.8 15 8 GC-FID 93.3 575-002 1-Chtoro-4-tritutoromethylberazene (parachtorobenzotrifluoride; Bilevel 75 14.41 15 8 GC-FID 93.3 575-002 Chtorobenzene Partial 75 14.41 15 8 GC-FID 93.3 575-002 Chtorobenzene Calculated 1000 20.02 GC-FID 575-001 575-002 Chtorobenzene (interple chetated) Calculated 100 50.C 13 60 8 GC-FID 75-001 Chtorotatene (methylchoride)	Carbon tetrachloride	Bilevel		10	25 C	14.1	30	8	GC-FID	98.3	575-001	
P-10102-2metry berzeer (monchronous) D12 13 15 6 GC-PD 91.6 97-401 P-10102-2metry berzeer (monchronous) Bilevel 25 0 11.8 15 8 GC-PD 91.6 97-501 P-10102-4rt/filuorometry/berzeer (parachlorobenzotr/filuoride; Bilevel 25 0 11.8 15 8 GC-PD 91.6 97-5002 P-10102-4rt/filuorometry/berzeer (parachlorobenzotr/filuoride; Bilevel 25 0 11.8 15 8 GC-PD 93.3 75-002 OXSOL 100 Chorobenzene Partial 75 14.41 15 8 GC-PD 93.6 75-002 Chorobenzene Partial 1000 20.02 GC-PD 875-001 60.8 GC-PD 875-001 Chorobenzene Gloval 1000 20.02 GC-PD 97.5 75-001 60.8 GC-PD 97.5 75-001 Chorobenzene (methylchoride) Calculated 10 50 12.8 15 8 GC-PD 97.5 75-001 <t< td=""><td>2-CELLOSOLVE acetate (2-ethoxyethyl acetate)</td><td>Calculated</td><td></td><td>100</td><td></td><td>12.1</td><td>15</td><td>0</td><td>GC-FID</td><td>95.4</td><td>575-002</td></t<>	2-CELLOSOLVE acetate (2-ethoxyethyl acetate)	Calculated		100		12.1	15	0	GC-FID	95.4	575-002	
Onlow Links (notoech conduct). Proceeding and the second	1-Chloro-2-methyl benzene (monochlorotoluene; OXSOL 10)	Bilevel		50 ‡		13	15	8	GC-FID	91.8 Q1	575-001	
OXSQL 100 Control Contro Control Control <	1-Chloro-4-(trifluoromethyl)benzene (parachlorobenzotrifluoride:	Bilevel		25 ◊		11.8	15	8	GC-FID	102	575-001	
1-Chiorx-4(trilluoromethyl)benzene (parachlorobenzotrifluoride; Bilevel 25 o 11.8 15 8 GC-FID 108 77-002 Chlorobenzene Partial 188 75 14.41 15 8 GC-FID 93.3 575-001 Chlorobenzene Partial 188 75 14.41 15 8 GC-FID 93.3 575-002 Chlorobenzene Calculated 1000 20.02 6 GC-FID 103 575-001 Chloroberhane (nethylchloride) Calculated 10 50C 13 60 8 GC-FID 575-001 Chloroberhane (nethylchloride) Calculated 24.6 GC-FID 575-001 575-001 Chloroberbane Bilevel 1374 50 ‡ 9.8 15 8 GC-FID 575-001 Cumene (sopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 105 575-001 Cyclohexane Bilevel 300 15.6 15 8 <td< td=""><td>OXSOL 100)</td><td></td><td></td><td>-</td><td></td><td></td><td>-</td><td></td><td></td><td></td><td></td></td<>	OXSOL 100)			-			-					
Chlorobenzene Partial 75 14.41 15 8 GC-FID 93.3 575-001 Chlorobenzene Partial 1838 75 14.41 15 8 GC-FID 87.6 575-002 Chlorobenzene Calculated 200 18.3 GC-FID 87.5 575-002 Chlorobenzene Calculated 1000 20.02 GC-FID 575-001 Chlorobenzene Bilevel 10 50 C 13 60 8 GC-FID 575-001 Chlorobenzene Dilevel 10 50 C 13 60 8 GC-FID 575-001 Chlorobenzene Calculated 12.4 GC-FID 575-001 575-001 Chlorobenzene Bilevel 50 12.8 15 8 GC-FID 105 575-001 Curnen (isopropyl benzene) Bilevel 300 15.6 15 8 GC-FID 105 575-001 Cyclohexane Bilevel 300 15.6	1-Chloro-4-(trifluoromethyl)benzene (parachlorobenzotrifluoride; OXSOL 100)	Bilevel		25 ◊		11.8	15	8	GC-FID	108	575-002	
Chorobanzene Partial 1838 75 14.41 15 8 GC-FID 87.6 87.5-002 Chlorobrance(hychloride) Calculated 1000 20.02 GC-FID 575-001 Chlorobrance(hychloride) Calculated 1000 20.02 GC-FID 575-001 Chlorobrance(hychloride) Calculated 100 50 C 13 60 8 GC-FID 575-001 Chlorobrance(hychloride) Calculated 10 50 C 13 60 8 GC-FID 575-001 Chlorobrance Bilevel 137 50 ‡ 9.8 15 8 GC-FID 575-002 Chlorobrance Gorpyt benzene) Bilevel 50 12.8 15 8 GC-FID 575-002 Cyclohexane Bilevel 50 12.8 15 8 GC-FID 105 575-002 Cyclohexane Bilevel 50 13.5 15 8 GC-FID 105 575-002 Cyclohexane	Chlorobenzene	Partial		75		14.41	15	8	GC-FID	93.3	575-001	
Chlorobromomethane Calculated 200 18.3 GC-FID 103 575-002 Chlorobrane (ethyl chloride) Calculated 1000 20.02 GC-FID 575-001 Chlorobrane (ethyl chloride) Calculated 10 50 C 13 60 8 GC-FID 575-001 Chlorobrane (ethyl chloride) Calculated 10 50 C 13 60 8 GC-FID 575-001 Chlorobrane (ethyl chloride) Calculated 50 ‡ 9.8 15 8 GC-FID 575-001 Chlorobrane (isopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 106 575-001 Curene (isopropyl benzene) Bilevel 300 15.6 15 8 GC-FID 106 575-001 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 108 575-001 Cyclohexane Calculated 50 13.5 GC-FID 108 575-001 Cyclohexane Calc	Chlorobenzene	Partial	1838	75		14.41	15	8	GC-FID	87.6	575-002	
Chloredhane (ethyl chloride) Calculated 100 20.02 CGC-FID 575-001 Chloroform Bilevel 10 50 C 13 60 8 GC-FID 575-001 Chloromethane (methylchloride) Calculated 24.6 GC-FID 575-001 575-001 o-Chlorostyrene Bilevel 1374 50 ‡ 9.8 15 8 GC-FID 572-001 Cumene (isopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 106 575-001 Cyclohexane Bilevel 50 12.8 15 8 GC-FID 106 575-001 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 108 575-001 Cyclohexane Calculated 50 13.5 GC-FID 105 575-001 Cyclohexane Calculated 50 13.5 GC-FID 105 575-001 Cyclohexane Calculated 300 15.4 GC-FID 106 </td <td>Chlorobromomethane</td> <td>Calculated</td> <td></td> <td>200</td> <td></td> <td>18.3</td> <td></td> <td></td> <td>GC-FID</td> <td>103</td> <td>575-002</td>	Chlorobromomethane	Calculated		200		18.3			GC-FID	103	575-002	
Chronotim Bilevel 10 50 C 13 60 8 GC-FID 97.3 57-001 Chromethane (methylchloride) Calculated 1374 50 ‡ 9.8 15 8 GC-FID 575-001 Choronethane (methylchloride) Calculated 12.4 GC-FID 97.3 575-001 Cumene (isopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 105 575-001 Cyclohexane Bilevel 50 12.8 15 8 GC-FID 105 575-002 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 105 575-001 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 105 575-001 Cyclohexane Calculated 50 13.5 GC-FID 105 575-001 Cyclohexane Calculated 300 15.4 GC-FID 102 575-001 Cyclohexene Calculated 3	Chloroethane (ethyl chloride)	Calculated		1000		20.02			GC-FID		575-001	
Chloribititatie (interrychlorible) Calculated Image: Calculated <thimage: calculated<="" th=""> Image: Calculated<!--</td--><td>Chloronorm Chloromethana (mathulahlarida)</td><td>Bilevel</td><td></td><td>10</td><td>50 C</td><td>13</td><td>60</td><td>8</td><td>GC-FID</td><td>97.3</td><td>575-001</td></thimage:>	Chloronorm Chloromethana (mathulahlarida)	Bilevel		10	50 C	13	60	8	GC-FID	97.3	575-001	
Onionational Control Dirty Dirty <thdirty< th=""> Dirty Dirty<td></td><td>Rilevel</td><td>1374</td><td>50 t</td><td></td><td>24.0</td><td>15</td><td>8</td><td>GC-FID</td><td>75.2</td><td>575-002</td></thdirty<>		Rilevel	1374	50 t		24.0	15	8	GC-FID	75.2	575-002	
Cumene (isopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 99.3 575-001 Cumene (isopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 106 575-002 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 106 575-002 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 105 575-002 Cyclohexanol Calculated 50 13.5 GC-FID 105 575-002 Cyclohexanol Calculated 50 13.5 GC-FID 106 575-002 Cyclohexanol Calculated 300 15.4 GC-FID 106 575-002 Cyclohexane Calculated 300 15.4 GC-FID 106 575-002 Cyclohexane Calculated 300 15.4 GC-FID 106 575-001 Cyclohexane Calculated 11.3 GC-FID 106 575-001	4-Chlorotoluene	Calculated	10/4	00 +		12.4	10		GC-FID	10.2	575-001	
Cumene (isopropyl benzene) Bilevel 50 12.8 15 8 GC-FID 106 575-002 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 105 575-001 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 105 575-002 Cyclohexanol Calculated 50 13.5 GC-FID 105 575-002 Cyclohexanol Calculated 50 13.5 GC-FID 105 575-002 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclopentane Calculated 300 15.4 GC-FID 106 575-001 Decamethylterporpytholuene) Calculated 300 15.4 GC-FID 106 575-001 Decamethylterporpytholuene) Calculated 11.3 GC-FID 575-001 575-001	Cumene (isopropyl benzene)	Bilevel		50		12.8	15	8	GC-FID	99.3	575-001	
Bilevel 300 15.6 15 8 GC-FID 105 575-001 Cyclohexane Bilevel 300 15.6 15 8 GC-FID 109 575-002 Cyclohexanol Calculated 50 13.5 C GC-FID 98 575-001 Cyclohexanol Calculated 50 13.5 C GC-FID 105 575-002 Cyclohexane Calculated 300 15.4 C GC-FID 105 575-001 Cyclohexene Calculated 300 15.4 C GC-FID 106 575-001 Cyclohexene Calculated 300 15.4 GC-FID 106 575-001 Cyclohexene Calculated 300 11.3 GC-FID 105 575-001 Cyclohexene Calculated 11.3 GC-FID 99 575-001 Decamethylopclopentasiloxane (D5) Partial 1891 5 † 5.666 15 8 GC-FID 90.3 <td< td=""><td>Cumene (isopropyl benzene)</td><td>Bilevel</td><td></td><td>50</td><td></td><td>12.8</td><td>15</td><td>8</td><td>GC-FID</td><td>106</td><td>575-002</td></td<>	Cumene (isopropyl benzene)	Bilevel		50		12.8	15	8	GC-FID	106	575-002	
Cyclohexane Bilevel 300 15.6 15 8 GC-FID 109 575-002 Cyclohexanol Calculated 50 13.5 GC-FID 105 575-002 Cyclohexanol Calculated 50 13.5 GC-FID 105 575-002 Cyclohexane Calculated 300 15.4 GC-FID 102 575-001 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclopentane Calculated 300 15.4 GC-FID 105 575-001 Decamethylopclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 9 575-001 Decamethylopclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 90 575-001 Decamethylopclopentasiloxane Calculated 7.36 GC-FID 97.3 <td>Cyclohexane</td> <td>Bilevel</td> <td></td> <td>300</td> <td></td> <td>15.6</td> <td>15</td> <td>8</td> <td>GC-FID</td> <td>105</td> <td>575-001</td>	Cyclohexane	Bilevel		300		15.6	15	8	GC-FID	105	575-001	
Cyclohexanol Calculated 50 13.5 GC-FID 98 575-001 Cyclohexanol Calculated 50 13.5 GC-FID 105 575-002 Cyclohexene Calculated 300 15.4 GC-FID 102 575-001 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclohexene Calculated 300 15.4 GC-FID 575-001 Cyclohexene Calculated 300 15.4 GC-FID 575-001 Cyclopentane Calculated 11.3 GC-FID 575-001 Decamethyleyclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 99 575-001 Decamethyleyclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 90 575-001 Decamethyleyclopentasiloxane Calculated 7.36 GC-FID 97.3 575-002 Decamethylettrasiloxane Partial	Cyclohexane	Bilevel		300		15.6	15	8	GC-FID	109	575-002	
Cyclonexanol Calculated 50 13.5 GC-FID 105 5/5-002 Cyclohexene Calculated 300 15.4 GC-FID 106 575-001 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclohexene Calculated 300 15.4 GC-FID 106 575-002 Cyclopentane Calculated 300 15.4 GC-FID 575-001 p-Cymene (4-isopropyltoluene) Calculated 11.3 GC-FID 575-001 Decamethyleyclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 99 575-001 Decamethyleyclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 99 575-001 Decamethyleyclopentasiloxane Calculated 7.36 GC-FID 97.3 575-002 Decamethyley alcohol Calculated 9.6 GC-FID 97.3 575-002 Decamethyletrasiloxane	Cyclohexanol	Calculated		50		13.5			GC-FID	98	575-001	
Cyclonexene Calculated 300 15.4 GC-FID 102 575-001 Cyclonexene Calculated 300 15.4 GC-FID 106 575-002 Cyclonexane Calculated 11.3 GC-FID 575-001 575-001 p-Cymene (4-isopropyltoluene) Calculated 11.3 GC-FID 575-001 Decamethylcyclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 99 575-001 Decamethylcyclopentasiloxane Calculated 7.36 GC-FID 99 575-001 Decamethylcyclopentasiloxane Calculated 7.36 GC-FID 103 575-002 Decamethylcyclopentasiloxane Calculated 9.6 GC-FID 97.3 575-002 Decamethylcyclopentasiloxane Calculated 9.6 GC-FID 94.3 575-002 Decamethylcyclopentasiloxane Partial 1893 2 13.8 30 4 GC-FID 94.3 575-002 Diacetone alcohol Cal		Calculated		50		13.5			GC-FID	105	575-002	
Optimizerie Calculated 300 13.4 Clock of the stress	Cyclohexene	Calculated		300		15.4			GC-FID	102	575-001	
Opcoment Calculated 10.0 Control Contro Control Control <t< td=""><td></td><td>Calculated</td><td></td><td>500</td><td></td><td>16.8</td><td></td><td></td><td>GC-FID</td><td>100</td><td>575-001</td></t<>		Calculated		500		16.8			GC-FID	100	575-001	
Decamethylcyclopentasiloxane (D5) Partial 1891 5 † 5.66 15 8 GC-FID 99 575-001 Decamethylcyclopentasiloxane Calculated 7.36 GC-FID 575-001 575-001 Decamethyltetrasiloxane Partial 12.2 GC-FID 103 575-002 n-Decane Partial 9.6 GC-FID 97.3 575-002 1-Decarlo (decyl alcohol (1-decarol) Calculated 9.6 GC-FID 97.3 575-002 Decyl alcohol (1-decarol) Calculated 9.6 GC-FID 94.3 575-002 Destflurane Partial 1893 2 13.8 30 4 GC-FID 94.3 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 Diacetone alcohol Calculated 1 ppb 12.6 GC-FID 91.3 575-002 Dibronochloromethane Calculated 1 ppb 12.6 GC-FID 92.9 575-001 Dibronochlor	p-Cymene (4-isopropyltoluene)	Calculated				11.3			GC-FID		575-001	
Decamethyltetrasiloxane Calculated 7.36 GC-FID 575-001 n-Decane Partial 12.2 GC-FID 103 575-002 1-Decanol (decyl alcohol) Calculated 9.6 GC-FID 97.3 575-002 Decyl alcohol (1-decanol) Calculated 9.6 GC-FID 97.3 575-002 Desflurane Partial 1893 2 13.8 30 4 GC-FID 94.3 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 Dibromo-3-chloropropane Calculated 19.b 12.6 GC-FID 92.9 575-002 Dibromochloromethane Calculated 15.6 GC-FID 92.9 575-001 L2-Dibromothoromethane Calculated 15.6 GC-FID 97.01 101.3 575-001	Decamethylcyclopentasiloxane (D5)	Partial	1891	5†		5.66	15	8	GC-FID	99	575-001	
Partial 12.2 GC-FID 103 575-002 1-Decanol (decyl alcohol) Calculated 9.6 GC-FID 97.3 575-002 Decyl alcohol (1-decanol) Calculated 9.6 GC-FID 97.3 575-002 Desflurane Partial 1893 2 13.8 30 4 GC-FID 94.3 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 L2-Dibromo-3-chloropropane Calculated 50 12.6 GC-FID 92.9 575-002 Dibromochloromethane Calculated 19.b 12.6 GC-FID 92.9 575-002 Dibromochloromethane Calculated 19.b 12.6 GC-FID 10.13 575-002 Dibromochloromethane Calculated 15.6 GC-FID 10.13 575-001	Decamethyltetrasiloxane	Calculated				7.36			GC-FID		575-001	
1-Decanol (decyl alcohol) Calculated 9.6 GC-FID 97.3 575-002 Decyl alcohol (1-decanol) Calculated 9.6 GC-FID 97.3 575-002 Desflurane Partial 1893 2 13.8 30 4 GC-FID 94.3 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 Diacetone alcohol Calculated 12.6 GC-FID 92.9 575-002 Dibromochloromethane Calculated 12.6 GC-FID 10.1 575-002 Dibromochloromethane Calculated 15.6 GC-FID 92.9 575-001 12-Dibromochloromethane (ethylene dibromide) Calculated 20 30 15.3 GC-FID 92.3 575-001	n-Decane	Partial				12.2			GC-FID	103	575-002	
Declaration (1-decanol) Calculated 9.6 GC-FID 97.3 575-002 Desflurane Partial 1893 2 13.8 30 4 GC-FID 94.3 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 J.2-Dibromo-3-chloropropane Calculated 1ppb 12.6 GC-FID 10.1 575-002 Dibromochloromethane Calculated 1pb 12.6 GC-FID 10.1 575-002 1_2-Dibromochloromethane Calculated 1pb 12.6 GC-FID 10.1 575-001 1_2-Dibromochloromethane Calculated 20 30 15.3 GC-FID 92.3 575-001	1-Decanol (decyl alcohol)	Calculated				9.6			GC-FID	97.3	575-002	
Destination Partial 1893 2 13.8 30 4 GC-HD 94.3 575-002 Diacetone alcohol Calculated 50 12.4 GC-FID 92.9 575-002 1,2-Dibromo-3-chloropropane Calculated 50 12.6 GC-FID 10.3 575-002 Dibromochloromethane Calculated 15.6 GC-FID 10.1 575-001 1,2-Dibromoethane (ethylene dibromide) Calculated 20 30 15.3 GC-FID 92.3 575-001	Decyl alcohol (1-decanol)	Calculated	1000		0	9.6	00	4	GC-FID	97.3	5/5-002	
Diabonic activitie Calculated 30 12.4 GC-FID 92.9 3/3-002 1,2-Dibromo-3-chloropropane Calculated 1 ppb 12.6 GC-FID 10.3 \$75-002 Dibromochloromethane Calculated 1 ppb 15.6 GC-FID 10.3 \$75-001 1_2-Dibromoethane (ethylene dibromide) Calculated 20 30 15.3 GC-FID 92.3 \$75-001	Desnurane	Partial	1893	50	2	13.8	30	4	GC-FID	94.3	575-002 575-002	
Calculated Ppo 12.0 Od FD 10.0 0.000 Z Dibromochloromethane Calculated 15.6 GC-FID 575-001 12-Dibromochloromethane (ethylene dibromide) Calculated 20 30 15.3 GC-FID 92.3 1575-001	1 2-Dibromo-3-chloronronane	Calculated		00 1 nnh		12.4			GC-FID	92.9 101 २	575-002	
1.2-Dibromoethane (ethylene dibromide) Calculated 20 30 15.3 GC-FID 92.3 575-001	Dibromochloromethane	Calculated		1 hhn		15.6			GC-FID	101.0	575-001	
	1,2-Dibromoethane (ethylene dibromide)	Calculated		20	30	15.3			GC-FID	92.3	575-001	

See page 212 for abbreviations.

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			OSHA	PELs 🛆		Samplir	ng Time					
	Validation	Research	TWA	CLG/STEL	Sampling Rate	Min	Мах	Analytical	DE %	SKC VOC	Chek 575	
Chemical Hazard	Level	Report	(ppm)	(ppm)	(ml/min)	(min)	(hrs)	Method	§	Cat. No.		
1,2-Dibromoethane (ethylene dibromide)	Calculated		20	30	15.3			GC-FID	99.4	575-002		
o-Dichlorobenzene (1 2-dichlorobenzene)	Partial			50 C	10.7	15	8	GC-FID GC-FID	79.2	575-001		-
o-Dichlorobenzene (1,2-dichlorobenzene)	Partial	1875		50	12.5	15	8	GC-FID	77.1	575-002		_
m-Dichlorobenzene (1,3-dichlorobenzene)	Calculated				12.95			GC-FID	91.8	575-001		
m-Dichlorobenzene (1,3-dichlorobenzene)	Calculated				12.95			GC-FID	92.7	575-002		
p-Dichlorobenzene (1,4-dichlorobenzene)	Calculated		75		12.95			GC-FID	91.1	575-001		
p-Dichlorobenzene (1,4-dichlorobenzene)	Calculated		/5		12.95			GC-FID GC-FID	94.7	575-002		
1.3-Dichlorobenzene (m-dichlorobenzene)	Calculated				12.95			GC-FID	92.7	575-002		
1,2-Dichlorobenzene (o-dichlorobenzene)	Partial			50 C	12.5	15	8	GC-FID	79.2	575-001		
1,2-Dichlorobenzene (o-dichlorobenzene)	Partial	1875		50 C	12.5	15	8	GC-FID	77.1	575-002		
1,4-Dichlorobenzene (p-dichlorobenzene)	Calculated		75		12.95			GC-FID	91.1	575-001		
1,4-Dichlorobenzene (p-dichlorobenzene)	Calculated		/5		12.95			GC-FID	94.7	575-002		_
Dichlorodifluoromethane (Freon 12)	Calculated		1000		18.6			GC-FID		575-001		_
1,1-Dichloroethane	Calculated		1000		16.85			GC-FID		575-001		
1,2-Dichloroethane (ethylene dichloride)	Bilevel		50	100	14.2	60	8	GC-FID	95.8	575-001		-
1,2-Dichloroethene (1,2-dichloroethylene)	Full				14.8	15	8	GC-FID	97.1	575-001		
1,1-Dichloroethene (vinylidene chloride)	Bilevel		1	15.0	12.3	60	8	GC-FID	95.2	575-001		
Dichloroethyl ether	Calculated		5 ‡	15 C	12.5	45	0	GC-FID	07.1	575-001		
1,2-Dichloroethylene (1,2-dichloroethene)	Full		200		14.8	15	8	GC-FID GC-FID	97.1	575-001		_
Dichloromethane (methylene chloride)	Full	1323	25	125	14.7	240	8π	GC-FID	96	575-001		
Dichloromethane (methylene chloride)	Full	1323	25	125	16	15	4	GC-FID	96	575-001		
1,2-Dichloropropane (propylene dichloride)	Bilevel		75		14.3	15	8	GC-FID	97.7	575-001		
1,1-Dichloropropene	Calculated				15.6		-	GC-FID		575-001		
cis-1,3-Dichloropropene	Partial	1886	1‡		13.6	15	8	GC-FID	101	575-002		
1 2-Dichlorotetrafluoroethane (Freon 114)	Calculated	1000	1000		14.4	15	0	GC-FID	99.4	575-002		
Dicyclopentadiene	Calculated		5 ±		11.8			GC-FID		575-001		
Diethyl ether (ethyl ether)	Calculated		400		16.4			GC-FID		575-001		
Diethyl ketone (3-pentanone)	Calculated		200 ‡		14.8			GC-FID	83.9	575-001		
Diethyl ketone (3-pentanone)	Calculated		200 ‡		14.8			GC-FID	100.3	575-002		
Diethylene glycol dimethyl ether (2-methoxyethyl ether)	Calculated				11.5			GC-FID		575-002		
Diethylene glycol monoethyl ether	Calculated				9.85			GC-FID		575-002		
Diethylene glycol monoethyl ether acetate	Calculated				9.88			GC-FID		575-002		
Diethylene glycol monomethyl ether (2-[2-methoxyethoxy] ethanol)	Calculated				11.3			GC-FID		575-002		
Diisobutyl ketone (DIBK), (isovalerone)	Bilevel	1305	50		10.3	30	8	GC-FID	98.3	575-002		
1,2-Dimethoxyethane (ethylene glycol dimethyl ether)	Calculated		1000		14.7			GC-FID		575-001	or 575-002	
Dimethol adipate	Calculated		1000		10.73			GC-FID GC-FID		575-001		_
Dimethyl disulfide	Calculated				15.4			GC-FID		575-001		_
2,5-Dimethyl hexane	Calculated				11.86			GC-FID		575-001		
2,2-Dimethyl methane	Calculated				21.7			GC-FID		575-001		
Dimethyl pentanedioate	Calculated				10.8			GC-FID		575-001		
Dimethyl sulfavida	Calculated				19			GC-FID		575-001		
N N-Dimethylaniline	Calculated		5		10.5			GC-FID		575-002		
2,2-Dimethylbutane (neohexane)	Calculated		0		14.2			GC-FID		575-001		
trans-1,2-Dimethylcyclohexane	Calculated				12.4			GC-FID	106.1	575-001		
N,N-Dimethylformamide (DMF)	Calculated		10		16.4			GC-FID	87.2	575-002		
2,3-Dimethylpentane	Calculated				12.8			GC-FID		575-001		
1,4-Dioxane	Calculated		100	-	15.8			GC-FID	91.4	575-002		_
Dipropyl ketone (4-hentanone)	Calculated		50 t		12.1			GC-FID	85.3	575-001		
Dipropyl ketone (4-heptanone)	Calculated		50 ‡		12.1			GC-FID	112	575-002		
Dipropylene glycol methyl ether	Calculated		100		10.8			GC-FID	84.3	575-002		
Dodecamethylcyclohexasiloxane	Calculated				6.75			GC-FID	10	575-001		
n-Dodecanol (lauryl alcohol)	Calculated				8.7			GC-FID	107.5	575-001		
1-Dodecene	Calculated				0./ 0.20			GC-FID	103	575-002		
1-Dodecyl alcohol (lauryl alcohol)	Calculated				8.7			GC-FID	107.5	575-001		
1-Dodecyl alcohol (lauryl alcohol)	Calculated				8.7			GC-FID	103	575-002		
Dodecyl methacrylate	Calculated				7.6			GC-FID		575-002		
Enflurane (ethrane)	Partial	1893		2	13.8	30	4	GC-FID	101	575-002		
Epichlorohydrin	Calculated		5		16.4			GC-FID	88.2	575-002		
2,3-Epoxypropyl methachylate (glycidul)	Calculated		50	1	11.8			GC-FID		575-002		
2.0 Lpoxypropyr methacrylate (gryoldyr methacrylate)	Calculated	L		L	11.40		L		L	515-00Z		

See page 212 for abbreviations.

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			OSHA	OSHA PELs A		Sampling Time					
	Validation	Research	TWA	CLG/STEL	Sampling Rate	Min	Max	Analytical	DE %	SKC VOC Chek 575	
Chemical Hazard	Level	Report	(ppm)	(ppm)	(ml/min)	(min)	(hrs)	Method	<u>§</u>	Cat. No.	
2-Ethowethanol	Calculated	1870	200		20.3	15	8	GC-FID	100.8	575-002	
2-Ethoxyethanol	Calculated		200		14.4			GC-FID	111.2	575-002	
2-Ethoxyethyl acetate (2-CELLOSOLVE acetate)	Calculated		100		12.1			GC-FID	95.4	575-002	
Ethrane (enflurane)	Partial	1893		2	13.8	30	4	GC-FID	101	575-002	
Ethyl acetate	Partial	1894	400		13.1	30	8	GC-FID	92.8	575-001	
Ethyl acetate	Partial	1894	400		14.1	30	8	GC-FID	100	575-002	
Ethyl acrylate	Bilevel	1070	5		13.7	15	8	GC-FID	94.2	575-002	
Ethyl alcohol (ethanol)	Partial	18/6	1000		20.3	15	8	GC-FID	100	575-002	
Ethyl benzene	Bilevel		100		12.9	15	6	GC-FID	100	575-001	
Ethyl benzene	OSHA 1002		100		13.83	15	8	GC-FID	99.1	575-002	
Ethyl bromide (bromoethane)	Calculated		200	250 #	18.5			GC-FID	00.1	575-001	
Ethyl butyl ketone (3-heptanone)	Calculated		50		12.2			GC-FID	87.9	575-001	
Ethyl butyl ketone (3-heptanone)	Calculated		50		12.2			GC-FID	103.4	575-002	
Ethyl chloride (chloroethane)	Calculated		1000		20.2			GC-FID		575-001	
Ethyl cyanide	Calculated				18.61			GC-FID		575-001	
Ethyl ether	Calculated		400		16.4			GC-FID		575-001	
Etnyl formate	Calculated		100		17.8			GC-FID	00	575-001	
2-Ethyl nethachylate	Full		100		9.0	15	8	GC-FID	84.7	575-002	
Ethyl methacrylate	Full		100		13.1	15	8	GC-FID	104	575-002	
Ethyl propionate	Calculated		100		14	10	0	GC-FID	104	575-001	
Ethyl tert-butyl ether (tert-butyl ethyl ether)	Bilevel	1356			13.1	15	8	GC-FID	101	575-001	
Ethylene dibromide (1,2-dibromoethane)	Calculated		20	30	15.3			GC-FID	92.3	575-001	
Ethylene dibromide (1,2-dibromoethane)	Calculated		20	30	15.3			GC-FID	99.4	575-002	
Ethylene dichloride (1,2-dichloroethane)	Bilevel		50	100 C	14.2	60	8	GC-FID	95.8	575-001	
Ethylene glycol	Calculated			100 mg/m ³ C	17.44			GC-FID		575-002	
Ethylene glycol diethyl ether	Calculated				12.27			GC-FID		575-002	
Ethylene glycol dimethyl ether acetate (butyl CELLOSOLVE acetate)	Calculated		5		14.7			GC-FID		575-001 OF 575-002	
Ethylene glycol monobexyl ether	Calculated		5		10.4			GC-FID		575-001	
Ethylene glycol monomethyl ether acetate (methyl CELLOSOLVE	Calculated		25		12.9			GC-FID	92.4	575-002	
acetate)											
Ethylene oxide	Full	1543	1	5 EL	21.2	15	8	GC-ECD	102	575-005	
2-Ethylhexanol	Calculated				10.93			GC-FID	93.7	575-002	
2-Ethyltoluene	Calculated				12.1			GC-FID	106	575-002	
3-Ethyltoluene	Calculated				12.1				01	575-002	
Freon 11 (trichlorofluoromethane)	Calculated				16.65			GC-FID	31	575-001	
Freon 113 (1.1.2-trichloro-1.2.2-trifluoroethane)	Calculated		1000	1250 #	14.1			GC-FID		575-001	
Freon 114 (1,2-dichlorotetrafluoroethane)	Calculated		1000		15.3			GC-FID		575-001	
Freon 12 (dichlorodifluoromethane)	Calculated		1000		18.6			GC-FID		575-001	
Glutaric acid dimethyl ester	Calculated				11.5			GC-FID		575-002	
Glycidol (2,3-epoxy-1-propanol)	Calculated		50		17.8			GC-FID		575-002	
Glycidyl methacrylate (2,3-epoxypropyl methacrylate)	Calculated	1000			11.45	15	0	GC-FID	00.0	575-002	
n-Heptone	Full Bilovol	1893	500		14.1	15	8	GC-FID	105	575-002	
n-Hentane	Bilevel		500		13.9	15	8	GC-FID	103	575-002	
4-Heptanone (dipropyl ketone)	Calculated		50 ±		12.1	10	0	GC-FID	85.3	575-001	
4-Heptanone (dipropyl ketone)	Calculated		50 ‡		12.1			GC-FID	112.2	575-002	
3-Heptanone (ethyl butyl ketone)	Calculated		50		12.2			GC-FID	87.9	575-001	
3-Heptanone (ethyl butyl ketone)	Calculated		50		12.2			GC-FID	103.4	575-002	
2-Heptanone (methyl n-amyl ketone)	Calculated		100		12.2			GC-FID	99.8	575-002	
1-Heptene	Calculated		0.00 +		13.1			GC-FID		575-001	
Hexachloropthane	Calculated		1		10.5			GC-FID		575-001	
Hexadecane	Calculated		1		77			GC-FID		575-001	
Hexamethyldisiloxane (L2)	Partial	1892	200 †		9.98	15	8	GC-FID	102.9	575-001	
Hexamethylene dibromide (1,6-dibromohexane)	Calculated				10.7			GC-FID		575-001	
n-Hexane	Bilevel		500		14.3	15	8	GC-FID	100	575-001	
n-Hexane	Bilevel		500		14.3	15	8	GC-FID	112	575-002	
Hexanol (hexyl alcohol)	Calculated	40=-			12.64		-	GC-FID	92.9	575-002	
2-Hexanone (methyl butyl ketone MBK)	Partial	1873	100		14.3	15	8	GC-FID	104	575-002	
Levene Hexone (methyl isobutyl ketone [MIRK1)	Bilevel	1304	100		14.5	30	8	GC-FID	94.6	575-001	
Hexone (methyl isobutyl ketone [MIBK])	OSHA 1004	1004	100		13.62	00	8	GC-FID	92.9	575-002	
sec-Hexyl acetate	Calculated		50		11.1		<u> </u>	GC-FID	02.0	575-002	
Hexyl alcohol (hexanol)	Calculated				12.64			GC-FID	92.9	575-002	
Hexylene	Calculated				14.5			GC-FID		575-001	

See page 212 for abbreviations.

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			OSHA	PELs 🛆		Samplir	ng Time			
Chemical Hazard	Validation Level	Research Report	TWA (ppm)	CLG/STEL (ppm)	Sampling Rate (ml/min)	Min (min)	Max (hrs)	Analytical Method	DE %	SKC VOC Chek 575 Cat. No.
Hexylene glycol	Calculated			25 C	11.81			GC-FID	-	575-002
lodomethane (methyl iodide)	Calculated				18.7			GC-FID		575-001
Isoamyl acetate	Calculated		100		11.9			GC-FID	91.9	575-001
Isoamyl acetate	Calculated		100	105 1	11.9			GC-FID	108	575-002
Isoamyl alcohol	Calculated		100	125 #	13.9			GC-FID	106	575-002
			150		12.0	15	Q	GC-FID	00.1	575-002
Isobutyl acrylate	Calculated		150		12.2	15	0	GC-FID	33.1	575-002
Isobutyl alcohol	Calculated		100		15.6			GC-FID	93	575-001
Isobutyl alcohol	Calculated		100		15.6			GC-FID	100	575-002
Isoflurane	Full	1893			13.2	15	8	GC-FID	96	575-002
Isooctyl alcohol	Calculated		50 ‡		10.9			GC-FID		575-002
Isopentane (2-methyl butane)	Calculated		1000	610 #	15.8			GC-FID		575-001
Isophorone	Calculated	1000	25	500 #	11.3	15	0	GC-FID	102 5	575-002
Isopropyl acetate	Calculated	1029	250	500 #	10.42	15	0	GC-FID	88.5	575-002
Isopropyl acetate	Calculated		250		14.2			GC-FID	101	575-002
Isopropyl alcohol (isopropanol)	Partial	1839	400	500 #	18.42	15	8	GC-FID	103.5	575-002
Isopropyl benzene (cumene)	Bilevel		50		12.8	15	8	GC-FID	99.3	575-001
Isopropyl benzene (cumene)	Bilevel		50		12.8	15	8	GC-FID	106	575-002
Isopropyl ether	Calculated		500		13.4			GC-FID		575-001
Isopropyl glycidyl ether	Calculated		50	50 #	12.8			GC-FID		575-001
4-Isopropyltoluene (p-cymene)	Calculated	1000	50		11.3		0	GC-FID	00.0	575-001
Isovalerone (diisobutyi ketone [DIBK])	Bilevel	1308	50		10.3	30	8	GC-FID	98.3	575-002
Lauryl alcohol (1-dodecanol)	Calculated				8.7			GC-FID	107.5	575-002
	Calculated				11 1			GC-FID	102	575-002
Mesityl oxide	Calculated		25		13.7			GC-FID	TOL	575-001
Mesitylene (1,3,5-trimethylbenzene)	Calculated		25 ¶		12.1			GC-FID	93.6	575-001
Mesitylene (1,3,5-trimethylbenzene)	Calculated		25 ¶		12.1			GC-FID	96	575-002
Methanol (methyl alcohol)	Partial	1895	200	250	1.2	15	8	GC-FID	101	575-007
2-Methoxy-1-propyl acetate	Calculated				12.1			GC-FID		575-002
1-Methoxy-2-propanol (propylene glycol monomethyl ether)	Calculated		100 ‡	150 #	14.7			GC-FID	82.9	575-001
1-Methoxy-2-propanol (propylene glycol monomethyl ether)	Calculated		100 ‡	150 #	14.7			GC-FID	100	5/5-002
acetate)	Calculated				12.2			GC-FID	100	575-001
1-Methoxy-2-propyl acetate (propylene glycol monomethyl ether	Calculated				12.1			GC-FID	103	575-002
acetate)										
2-Methoxyethanol (methyl CELLOSOLVE)	Calculated		0.1		16.1			GC-FID	94.7	575-001
2-Methoxyethanol (methyl CELLOSOLVE)	Calculated		0.1		16.1			GC-FID	91.1	575-002
2-[2-Methoxyethoxy] ethanol (diethylene glycol monomethyl ether)	Calculated				11.3			GC-FID		575-002
Methoxyfurane	Calculated			2	13.3			GC-FID	95.7	575-002
Methyl acetate	Calculated		200	250 #	17.8			GC-FID	00.1	575-002
Methyl acrylate	Full		10		15.7	15	8	GC-FID	94.3	575-002
Methyl alcohol (Methanol)	Partial	1895	200	250	1.2	15	8	GC-FID	101	575-007
Methyl amyl alcohol (methyl isobutyl carbinol)	Calculated		25		12.8			GC-FID		575-002
Methyl bromide (bromomethane)	Calculated			20C	22.1			GC-FID		575-002
2-Methyl butane (isopentane)	Calculated	1070	1000	610 #	15.8	45	0	GC-FID	101	575-001
Methyl butyl ketone (MBK), (2-nexanone)	Coloulated	18/3	100		14.3	15	8	GC-FID	104	575-002
Methyl CELLOSOLVE (2-methoxyethanol)	Calculated		0.1		16.1			GC-FID	94.7	575-002
Methyl CELLOSOLVE (2 methodyourdator)	Calculated		25		12.9			GC-FID	92.4	575-002
acetate)										
Methyl chloroform (1,1,1-trichloroethane)	Bilevel		350		14.1	15	8	GC-FID	99.9	575-001
Methyl cyclohexane	Bilevel		500		14.2	15	8	GC-FID	106	575-001
Methyl ethyl ketone (MEK), (2-butanone)	Bilevel	1306	200		17.1	15	12	GC-FID	100	575-002
Methyl ethyl ketone (MEK), (2-butanone)	OSHA 1004		200		16.88		8	GC-FID	92.3	575-002
Methyl iormale	Calculated		100		20.58			GC-FID		575-001
Methyl isoamyl ketone	Calculated		50		12.3			GC-FID		575-002
Methyl isobutyl carbinol (methyl amvl alcohol)	Calculated		25		12.8			GC-FID		575-002
Methyl isobutyl ketone (MIBK), (hexone)	Bilevel	1304	100		13.5	30	8	GC-FID	94.6	575-002
Methyl isobutyl ketone (MIBK), (hexone)	OSHA 1004		100		13.62		8	GC-FID	92.9	575-002
Methyl isopropyl ketone	Calculated				14.8			GC-FID		575-002
Methyl isothiocyanate	Calculated				17.36			GC-FID		575-001
Methyl methacrylate (MMA)	Bilevel	1308	100		13.1	7.5	8	GC-FID	100.5	575-002
Methyl n-amyl ketone (2-heptanone)	Calculated		100		12.2			GC-FID	99.8	575-002
Z-mernyi pentane	Calculated		000		14.1			GC-FID	00.0	575-001
Internyi propyi ketone (2-pentanone)	Calculated		200	L	14.0		L	GO-FID	92.0	373-002

See page 212 for abbreviations.

			OSHA	PELs 🛆		Samplin	ng Time			
	Validation	Research	TWA	CLG/STEL	Sampling Rate	Min	Max	Analytical	DE %	SKC VOC Chek 575
Chemical Hazard	Level	Report	(ppm)	(ppm)	(ml/min)	(min)	(hrs)	Method	§	Cat. No.
S-Methyl styrene (vinyl toluene)	Calculated		100		12.20			GC-FID		575-002
Methyl t-butyl ether (MTBE)	Full	1352			13.6	8.5	8	GC-FID	97.4	575-001
Methyl tert-amyl ether (tert-amyl methyl ether)	Bilevel	1355			13.1	30	8	GC-FID	99	575-001
n-Methyl-2-pyrrolidinone	Calculated		51		13.97			GC-FID		575-001
5-Methyl-3-heptanone	Calculated		25		11.4			GC-FID	87.5	575-001
5-Methyl-3-neptanone	Calculated		25		11.4			GC-FID	110.7	575-002
Methylal (diffectioxymethalie)	Calculated		1000		24.6			GC-FID		575-002
1-Methylcyclohexanol	Full		100		12.4	15	8	GC-FID	94.7	575-001
1-Methylcyclohexanol	Full		100		12.4	15	8	GC-FID	108	575-002
Methylcyclopentane	Calculated	1000		105	14.9			GC-FID		575-001
Methylene chloride (dichloromethane)	Full	1323	25	125	16	15 240	4 8 m	GC-FID	96	575-001
3-Methylhexane	Calculated	1020	25	125	12.8	240	0 /	GC-FID	30	575-001
alpha-Methylstyrene	Bilevel	1359		100 C	12.6	15	12	GC-FID	95.7	575-002
Mineral spirits	Calculated		500		10.95			GC-FID		575-001
Monochlorotoluene (1-chloro-2-methylbenzene; OXSOL 10)	Bilevel		50 ‡		13	15	8	GC-FID	91.8	575-001
Monochlorotoluene (1-chloro-2-methylbenzene; OXSOL 10)	Bilevel		50 ‡		13	15	8	GC-FID	91	575-002
Neonexane (2,2-dimethyibutane)	Calculated				14.2			GC-FID GC-FID		575-001
Nonane	Bilevel				10.6	15	8	GC-FID	103	575-001
Nonyl alcohol	Calculated				10.2			GC-FID	96.8	575-002
Octadecane	Calculated				7.1			GC-FID		575-001
Octamethylcyclotetrasiloxane (D4)	Partial	1890	10 🔻		6.32	15	8	GC-FID	97.2	575-001
Octamethyltrisiloxane (L3)	Partial	1902	200 †		8.47	15	8	GC-FID	98.3	575-001
	Bilevel		500		12.7	15	8	GC-FID	110	575-001
Octanol (octyl alcohol)	Calculated		500		10.86	15	0	GC-FID	110	575-002
1-Octene	Calculated				11.99			GC-FID		575-001
Octyl alcohol (octanol)	Calculated				10.86			GC-FID		575-002
OXSOL 10 (monochlorotoluene [1-chloro-2-methyl benzene])	Bilevel		50 ‡		13	15	8	GC-FID	91.8	575-001
OXSOL 10 (monochlorotoluene [1-chloro-2-methyl benzene])	Bilevel		50 ‡		13	15	8	GC-FID	91	575-002
OXSOL 100)	Dilever		25 V		11.0	15			102	575-001
Parachlorobenzotrifluoride (1-chloro-4-[trifluoromethyl]benzene;	Bilevel		25 ◊		11.8	15	8	GC-FID	108	575-002
OXSOL 100)										
Pentadecane	Calculated	1011	1000		7.93	15	0	GC-FID	105.0	5/5-001
3-Pentanone (diethyl ketone)	Calculated	1311	200 ±		14.9	15	0	GC-FID	83.9	575-001
3-Pentanone (diethyl ketone)	Calculated		200 ‡		14.8			GC-FID	100.3	575-002
2-Pentanone (methyl propyl ketone)	Calculated		200		14.8			GC-FID	92.6	575-002
1-Pentene	Calculated				16.3			GC-FID		575-001
2-Pentyl acetate (sec-amyl acetate)	Calculated	1000	125	000.0	11.8	7.5	10	GC-FID	100.0	575-001
Perchloroethylene (tetrachloroethylene)		1080	100	200 C	13.1	7.5	8	GC-FID	95.4	575-001
Perfluoromethylcyclohexane	Partial		100	200 0	10.00	5	0	GC-FID	102	575-002
Phenol (carbolic acid)	Calculated		5	15.6 C	14.5			GC-FID		575-001 or 575-002
Phenyl ether (diphenyl oxide)	Calculated		1		10.4			GC-FID		575-001
Phenyl glycidyl ether	Calculated		10		11.6			GC-FID		575-001
4-menyicyclonexene	Partial	1840			11.53	15	8	GC-FID	108.6	575-001 or 575-002
Propane	Calculated	1040	1000		21.73	10	0	GC-FID	100.0	575-001
n-Propanol (propyl alcohol)	Calculated		200		17.7			GC-FID	87.3	575-001
n-Propanol (propyl alcohol)	Calculated		200		17.7			GC-FID	97.8	575-002
Propionitrile	Calculated		6‡		18.61			GC-FID		575-001
n-Propyl acetate	Calculated		200		14.1			GC-FID	87.5	575-002
Propyl alcohol (n-propanol)	Calculated		200		14.1			GC-FID	87.3	575-001
Propyl alcohol (n-propanol)	Calculated		200		17.7			GC-FID	97.8	575-002
Propyl bromide (1-bromopropane)	Full	1740	0.1		14.4	30	8	GC-FID	100	575-001
Propyl bromide (1-bromopropane)	Full	1740	0.1		14.7	30	8	GC-FID	107	575-002
n-Propylbenzene	Calculated		75		12.1	15		GC-FID	101	575-002
Propylene dichloride (1,2-dichloropropane) Propylene divcol monomethyl ether (1-methoxy-2-propanol)	Calculated		/5	150 #	14.3	15	8	GC-FID	97.7 82 Q	575-001
Propylene glycol monomethyl ether (1-methoxy-2-propanol)	Calculated		100 ±	150 #	14.7			GC-FID	100	575-002
Propylene glycol monomethyl ether acetate (1-methoxy-2-propyl	Calculated		· • • T		12.2			GC-FID	108	575-001
acetate)										
Propylene glycol monomethyl ether acetate (1-methoxy-2-propyl	Calculated				12.1			GC-FID	103	575-002
averare)										

•

See page 212 for abbreviations.

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			OSHA	PELs A		Samplin	ng Time			
	Validation	Besearch	TWA	CI G/STEL	Sampling Bate	Min	Max	Analytical	DF %	SKC VOC Chek 575
Chemical Hazard	Level	Report	(ppm)	(ppm)	(ml/min)	(min)	(hrs)	Method	§	Cat. No.
Propylene oxide	Calculated		100		19.9			GC-FID	98	575-001
Propylene oxide	Calculated		100		19.9			GC-FID	99.7	575-002
Pyridine	Calculated		5		16.6			GC-FID	88.2	575-002
Sevoflurane	Partial	1893		2	12.8	30	4		100	575-002
Solvent naphtha (petroleum) light aromatic	Calculated				11.61			GC-FID		575-001 or 575-002
Stoddard solvent	Calculated		500		10.95			GC-FID		575-001
Styrene	Full	1313	100	200 C	13.7	15	8	GC-FID	86.3	575-002
Styrene	OSHA 1014		100	200 C	13.55	15	8	GC-FID	96.7	575-006
1,1,1,2-Tetrachloroethane	Calculated				13.63			GC-FID	98.3	575-002
1,1,2,2-Tetrachloroethane	Bilevel		5		11.8	480 *	8	GC-FID	64.4 ∞	575-001
Tetrachloroethylene (perchloroethylene)	Full	1686	100	200 C	13.1	7.5	12	GC-FID	100.8	575-001
Tetrachloroethylene (perchloroethylene)	OSHA 1001		100	200 C	13.06	5	8	GC-FID	95.4	575-002
Tetradecane	Calculated				8.3			GC-FID		575-001
Tetrahydrofuran	Partial	1841	200		17.7	15	8	GC-FID	100.6	575-002
1,2,3,4-Tetramethylbenzene	Calculated				11.1			GC-FID		575-001
1,2,3,5-Tetramethylbenzene	Calculated				10.8			GC-FID		575-001
1,2,4,5-Tetramethylbenzene	Calculated				11.2			GC-FID	86.6	575-002
Toluene	Bilevel		200	300 C	14.5	15	8	GC-FID	97.9	575-001
Toluene	OSHA 111		200	300 C	14.89	10	8	GC-FID	97	575-002
1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113)	Calculated		1000	1250 #	14.1			GC-FID		575-001
1,2,3-Trichlorobenzene	Calculated				11.34			GC-FID		575-001
1,2,4-Trichlorobenzene	Calculated			5 C	11.4			GC-FID		575-001
1,1,2-Trichloroethane	Bilevel		10		12.5	15	8	GC-FID	96.7	575-001
1,1,1-Trichloroethane (methyl chloroform)	Bilevel		350		14.1	15	8	GC-FID	99.9	575-001
Trichloroethylene	Full		100	200 C	14.9	15	8	GC-FID	102	575-001
Trichloroethylene	OSHA 1001		100	200 C	14.24	5	8	GC-FID	97.5	575-002
Trichlorofluoromethane (Freon 11)	Calculated				16.65			GC-FID		575-001
1,2,3-Trichloropropane	Bilevel		50		11.9	15	8	GC-FID	98.1	575-001
Tridecane	Calculated				9.2			GC-FID		575-001
Trifluoromethyl benzene (benzotrifluoride; OXSOL 2000)	Bilevel		100		13.3	15	8	GC-FID	106	575-001
Trifluoromethyl benzene (benzotrifluoride; OXSOL 2000)	Bilevel		100		13.3	15	8	GC-FID	107	575-002
1,2,3-Trimethylbenzene	Calculated		25 ¶		12			GC-FID	91.1	575-001
1,2,3-Trimethylbenzene	Calculated		25 ¶		12	15	8	GC-FID	93.8	575-002
1,2,4-Trimethylbenzene	Partial		25 ¶		13.05			GC-FID	88.4	575-001
1,2,4-Trimethylbenzene	Partial	1837	25 ¶		13.05	15	8	GC-FID	88.9	575-002
1,3,5-Trimethylbenzene (mesitylene)	Calculated		25 ¶		12.1			GC-FID	93.6	575-001
1,3,5-Trimethylbenzene (mesitylene)	Calculated		25 ¶		12.1			GC-FID	96	575-002
2,2,4-Trimethylpentane	Calculated				11.89			GC-FID		575-001
n-Undecane	Calculated				9.62			GC-FID		575-001
Vinyl acetate	Full	1860		4 C #	16.4	30	8	GC-FID	92	575-002
Vinyl bromide	Calculated		LFC ‡		19.6			GC-FID		575-001
Vinyl chloride	Calculated		1		21.4			GC-FID		575-001
Vinyl toluene (methyl styrene)	Calculated		100		12.3			GC-FID		575-001
n-Vinyl-2-pyrrolidone	Calculated				13.9			GC-FID		575-001
4-Vinylcyclohexene	Calculated				12.4		-	GC-FID		575-001
Vinylidene chloride (1,1-dichloroethene)	Bilevel		1		12.3	60	8	GC-FID	95.2	575-001
m-Xylene	Bilevel		100	150 #	12.5	15	8	GC-FID	96.6	575-001
m-xylene	Bilevel		100	150 #	12.5	15	8	GC-FID	101	575-002
m-xyiene	USHA 1002		100	150 #	13.82		8	GC-FID	96.1	575-002
o-Xylene	Bilevel		100	150 #	11.9	15	8	GC-FID	91	575-000
0-Xylene	OSHA 1002		100	150 #	14.24		8	GC-FID	89.4	575-002
p-Xylene	Bilevel		100	150 #	12.8	15	8	GC-FID	95.6	575-001
p-Xylene	Bilevel		100	150 #	12.8	15	8	GC-FID	103	575-002
p-Xylene	OSHA 1002		100	150 #	13.94		8	GC-FID	95.3	575-002

Lower than the NIOSH-accepted guideline

NIOSH Short Term Exposure Limit (STEL) #

Depends on detector sensitivity 00

In-house exposure level †

NIOSH Recommended Exposure Limit (REL) Occidental Chemical corporate exposure limits ‡ Ó

≈

Valid for PEL samples greater than 4 hours duration. If more than 1000 ppm of other contaminants are present, reduce maximum sample time to 4 hours.

Agency standards for OSHA listings represent the OSHA PELs reported in 29 CFR 1910.1000 Part 1910, Section 1000. Δ

§ The values given for the desorption efficiency were obtained in SKC Inc. laboratories. See the online guide at www.skcinc.com for details on the desorption solvent used.

OSHA construction industry standards

ſ Valid for STEL samples up to 4 hours duration $\sqrt{}$

If more than 1000 ppm contaminants are present, reduce maximum sample time to 4 hours. OARS-WEEL TWA Level (Occupational Alliance for Risk Science - Workplace Environmental Exposure Levels) \mathbf{v}

EL Excursion Limit

LFC Lowest feasible concentration



See page 212 for abbreviations.

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Passive Samplers Halogenated Ethers/Ethylene Oxide

New design easier analysis

New design

easier analysis

Rack

Passive Samplers for Health Care Anesthetic Gases

From

Halogenated Ethers

Isoflurane and enflurane are used primarily in veterinary procedures. Isoflurane and halothane are being replaced by desflurane and sevoflurane, which have fewer health effects. Exposure to halogenated ethers can cause eye, skin, and respiratory tract irritation; eye damage; and other serious health effects.

SKC VOC Chek 575 Passive Samplers

- Validated sampling rates for each gas
- Lightweight and miniature size
- No pump or assembly required
- Easy sorbent transfer during analysis

	Pkg. of 5	Pkg. of 25
Description	Cat. No.	Cat. No.
Organic Vapors Passive Sampler contains 500 mg Anasorb 747	575-002 [†]	575-002A [†]
t Larger quantity packages are available. Contact SKC		

See details on pages 80-81.

Fron

BO

Sampling Anesthetic Gases OSHA Anesthetic Gases Guidelines for Workplace Exposures recommend air sampling for anesthetic gases every six months to measure worker exposures and to check control measure effectiveness. Personal sampling provides the best estimate of exposure level and is the preferred method for determining worker time-weighted average (TWA) exposure during anesthetic administration and in the post-anesthesia care unit (PACU). Go to www.osha.gov/dts/ osta/anestheticgases/index.html for detailed guidelines.

BO Sampling Peracetic Acid (PAA) -

an Alternate Sterilant

While no passive sampler is available for PAA, SKC offers treated sorbent tubes Cat. Nos. 226-193-UC and 226-199-UC (page 42) and coated filter Cat. No. 225-9030 (page 63) to be used in a sampling train for PAA and hydrogen peroxide. See page 42 for details.



Ethylene Oxide (EtO)

Ethylene oxide is used to sterilize medical equipment and supplies. Exposure health effects can range from respiratory irritation to cancer and reproductive/mutagenic effects.

SKC Ethylene Oxide Passive Sampler

- Uses same sorbent as active OSHA Method 1010
- Fully validated for 8-hour and 15-minute sampling
- Lightweight and miniature size
- No pump or assembly required

	Pkg. of 5	Pkg. of 25
Description	Cat. No.	Cat. No.
Ethylene Oxide Passive Sampler contains 500 mg Anasorb 747 treated		
with hydrobromic acid	575-005 [†]	575-005A [†]
t Larger swartity people are subjected CVC		

Larger quantity packages are available. Contact SKC

For passive samplers for Aldehydes, see page 84.

UME^X 100 Formaldehyde

UME^X100 Passive Sampler for Formaldehyde Occupational and Indoor Air Exposure Monitoring

Sampling Rates for Other Aldehydes

Compound	Sampling Rate (ml/min)
Formaldehyde	28.6
(full validation)	(velocity 5 to 100 cm/
	sec, 15 min to 24 hrs)
	20.4
	(velocity < 5 cm/sec,
	1 to 7 days)
Acetaldehyde	22.8 [‡]
Benzaldehyde	13.5 [‡]
Butyraldehyde	15.8 [‡]
Crotonaldehyde	9.71‡
Glutaraldehyde	14 [‡]
Hexanaldehyde	9.66 [‡]
Isovaleraldehyde	15.5 [‡]
Propionaldehyde	14 [‡]
Chloroacetaldehyde	19.4**
Decylaldehyde	10.4**
Heptanaldehyde	12.8**
Nonanaldehyde	11.6**
o-Phthaldehyde	12.83**
o-Tolualdehyde	12.7**
Valeraldehyde	15.4**

‡ Partial Validation

Calculated sampling rate; see online Passive Sampling Guide at www.skcinc.com

C Tech Tips

Q: Why is my lab finding higher background levels of formaldehyde on sampling media than those reported by SKC?

A: Page 3 of OSHA 1007 reports that storing samplers at elevated temperatures will cause DNPH to decompose. The decomposition product, 2,4-dinitroaniline, may be seen as formaldehyde by some labs if they are using a short/fast HPLC column with inadequate plate count. EPA TO-11A, Section 13.2.2 recommends an HPLC column efficiency of > 5000 theoretical plates. This column will allow for the effective separation of the formaldehyde peak from interfering peaks.

- Meets OSHA Method 1007 specifications
- Conforms to EU ISO 16000-4-2004
- Accuracy exceeds OSHA standards
- Highly selective 2,4-DNPH chemistry; easy analysis
- Documented formaldehyde uptake rates for 15-minute to 24-hour and 7-day samples
 - Sampling rates available for other aldehydes (see left)
- Samples low ppb levels of formaldehyde



Versatile UME^x 100 Sampler



Easy **personal formaldehyde sampling** (15 minutes to 8 hours) — OSHA 1007



Convenient **indoor air formaldehyde sampling** (24 hours or 7 days) with UME^X 100 and stand accessory

	# # #		
Description	Choines	Cat. No.	Price/Qty.
UME ^x 100 Passive Sampler for Formaldehyde*t	A01067	500-100	\$ 159.00/10
Suitable for sampling other aldehydes		500-100A	379.00/25
Treated Tape for QC - UME ^x 100* [†]		P20084	35.00/50
Accessory		Cat. No.	Price/Qty.
Stand for Area Sampling		690-302	\$ 8.00/ea

Limited shelf-life, single use only; do not reuse

Store at \leq 4 C (39.2 F)

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 Δ If sampling in an atmosphere containing formalin, see www.skcinc.com/instructions/1795.pdf.

UME^x 200 and UME^x 300

Nitrogen Dioxide/Sulfur Dioxide and Ammonia

UME^x200 Passive Sampler for NO₂ and SO₂

Occupational Sampling and Near-road/Urban Air Monitoring

- Same chemistry as active OSHA Method ID-182
- Accurately measures exposures to sulfur dioxide and/or nitrogen dioxide from 15 minutes to 24 hours
- 3-week sample storage at ambient temperature
- Documented sampling rate of 17.3 ml/min for NO₂ and 15.2 ml/min for SO₂





Cat. No. Price/Qty. Stand for Area Sampling 690-302 \$ 8.00/ea



Personal SO₂/NO₂ sampling (15 minutes to 8 hours)



Fenceline monitoring (up to 24 hours) with UME^X 200 and shelter



Occupational exposure monitoring in:

- Chemical plants
- Textile manufacturing
- Food industry
- Copper smelting
- Power plants
- Paper pulp mills
- Cement manufacturing
- Mines
- Welding

Standard of

- Construction
- Fertilizer production Explosives production

UME^x 200 is ideal for:

- Near-road monitoring
- Fenceline applications
- Urban air studies by community action groups

Good Practice

Store and prepare sampling media in solvent-free environments.

Limited shelf-life, single use only; do not reuse

UME^x300 Passive Sampler for Ammonia

Occupational Sampling and Fenceline/Odor Monitoring

- Chemistry similar to active OSHA Method **ID-188 and NIOSH 6016**
- Safe no glass or sulfuric acid liquid in the sampler
- No particulate interference
- Enhanced sensitivity with documented 39.92 ml/min uptake rate
- Detects ammonia to: • 2.4 ppm for a 15-minute sample
 - 0.075 ppm for an 8-hour sample
 - 0.025 ppm for a 24-hour sample

/*		
Description	Cat. No.	Price/Qty.
UME ^X 300 Passive Sampler for		
Ammonia*†	500-300	\$ 169.00/10
Treated Tape for QC - UME ^x 300* [†]	P20083	39.00/25
	2 -	
Accessory	Cat. No.	Price/Qty.
Stand for Area Sampling	690-302	\$ 8.00/ea

Limited shelf-life, single use only; do not reuse

Store at $\leq 4 C (39.2 F)$

SKC Asia www.skc-asia.com SKC Inc.



Personal ammonia sampling (15 minutes to 8 hours)



Fenceline monitoring (24 hours) with UMEX 300 and stand accessory

www.skcinc.com

UME^x 300 Sampler Applications Ammonia is one of the most

ΑΒΟΙ

commonly produced industrial chemicals in the U.S and is found in the following industries:

- Agricultural (fertilizers, poultry, swine, dairy)
- Janitorial services (cleaners)
- Food and beverage (r efrigeration and fermentation)
- Remediation
- Fuel
- Manufacturing (plastics,
- explosives, pesticides, textiles, dyes)

Passive Samplers Mercury/HCN



The 520 Elemental Mercury Passive Sampler only samples elemental mercury (Hg) in the vapor phase; it does not sample elemental Hg in the particulate phase or organic mercury compounds.



- Elemental Mercury Passive Sampler For OSHA Method ID-140
- Lowest cost per measurement available
 - Reusable capsule holder
 - Replaceable sorbent capsule
- Lightweight and easy to use; no pump needed
- No moisture or chlorine interferences
- Long-term sampling up to 120 hours
- Validated by OSHA ID-140
- High accuracy, sensitivity, and capacity
 Positive analysis of mercury
 - Removable sorbent capsule eliminates false high readings due to contamination of capsule holder

The SKC Elemental Mercury Passive Sampler measures worker exposure level as a time-weighted average (TWA) and permits positive analysis for mercury vapor. This economical and reliable passive sampler is designed for analysis by atomic absorption. The capsule holder can be cleaned and reused to reduce sampling costs.



Description	Cat. No.	Qty.			
Sorbent Capsules contain Anasorb C300* and include replacement foams and					
resealable bags	520-02A	10			
-	520-02C	30			
Reusable Capsule Holder	520-03	ea			
,	520-03A	5			
Anasorb C300 is equivalent to Carulite and Hydrar.					

Note: To sample low levels of mercury, use a sorbent tube.

Hydrogen Cyanide Passive Sampler For OSHA Method 1015

- More accurate than previous methods
- Lightweight miniature sampler for unobtrusive sampling
- Samples hydrogen cyanide from 0.44 to 20 ppm
 - Suitable for 15-minute and 8-hour samples
- Unique sampler design for direct transfer of sorbent into and out of the sampler
- Back view Erc
- Load and unload sorbent easily in the field
 Sorbent transfers directly to vial for solvent extraction



SKC offers the Hydrogen Cyanide (HCN) Passive Sampler specified in OSHA 1015. The HCN Passive Sampler operates at a sampling rate of 28.4 ml/min and provides accurate HCN exposure results. The sampler design allows field loading of sorbent from vial to sampler housing before sampling and field transfer of sorbent from sampler to vial after sampling. Samples are extracted with water and analyzed by ion chromatography/electrochemical detector (IC/ELCHM).

Description	Cat. No.	Qty.
Hydrogen Cyanide Passive Sampler includes empty sampler housing and sealed glass		
vial of 600-mg soda lime sorbent	590-400	5