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APPENDIX A

APPENDIX A

Calibration Curve of BTEX and MTBE

```
=====
Calibration Table
=====
```

Calib. Data Modified : 2/4/03 12:31:12 PM
 Calculate : External Standard
 Based on : Peak Area
 Rel. Reference Window : 5.000 %
 Abs. Reference Window : 0.000 min
 Rel. Non-ref. Window : 5.000 %
 Abs. Non-ref. Window : 0.000 min
 Uncalibrated Peaks : not reported
 Partial Calibration : Yes, identified peaks are recalibrated
 Correct All Ret. Times: No, only for identified peaks
 Curve Type : Linear
 Origin : Forced
 Weight : Equal
 Recalibration Settings:
 Average Response : Average all calibrations
 Average Retention Time: Floating Average New 75%
 Calibration Report Options :
 Printout of recalibrations within a sequence:
 Calibration Table after Recalibration
 Normal Report after Recalibration
 If the sequence is done with bracketing:
 Results of first cycle (ending previous bracket)

Signal 1: FID1 A,

RetTime	Lvl	Amount	Area	Amt/Area	Ref Grp	Name
[min]	Sig	[ug]				
6.575	1	9.90100	1.12959	8.76509	MTBE	
	2	47.61900	6.55217	7.26767		
	3	90.90900	14.05666	6.46733		
	4	230.77000	36.09025	6.39425		
8.945	1	4.97500	1.64325	3.02753	BENZENE	
	1	9.90100	2.27270	4.35649		
	2	47.61900	11.57973	4.11227		
	3	90.90900	23.97835	3.79130		
	4	230.77000	59.72768	3.86370		

RetTime [min]	Lvl Sig	Amount [ug]	Area	Amt/Area	Ref Grp	Name
12.671	1 5	4.97500	1.33320	3.73163	TOLUENE	
	1	9.90100	2.32718	4.25451		
	2	47.61900	12.21735	3.89765		
	3	90.90900	25.31466	3.59116		
	4	230.77000	64.29553	3.58921		
16.401	1 5	4.97500	1.40644	3.53730	ETHYLBENZENE	
	1	9.90100	2.38735	4.14727		
	2	47.61900	12.87799	3.69770		
	3	90.90900	25.95047	3.50317		
	4	230.77000	67.51861	3.41787		
16.696	1 5	4.97500	1.35628	3.66813	M-XYLENE	
	1	9.90100	2.53936	3.89901		
	2	47.61900	13.31071	3.57750		
	3	90.90900	26.75976	3.39723		
	4	230.77000	69.24318	3.33275		
17.550	1 5	4.97500	1.62161	3.06794	O-XYLENE	
	1	9.90100	2.74222	3.61057		
	2	47.61900	13.61374	3.49786		
	3	90.90900	27.38919	3.31916		
	4	230.77000	70.53236	3.27183		

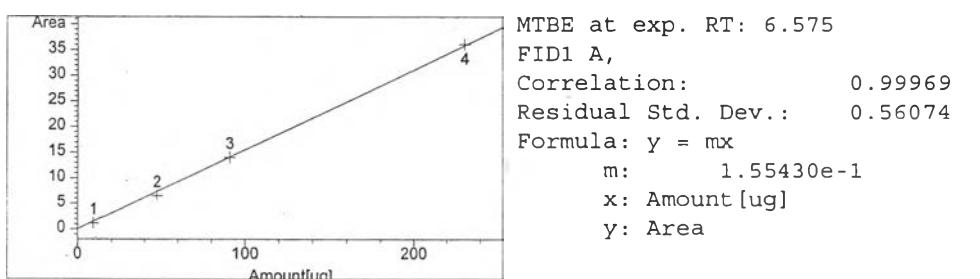
2 Warnings or Errors :

Warning : Overlapping peak time windows at 16.401 min, signal 1
 Warning : Overlapping peak time windows at 16.696 min, signal 1

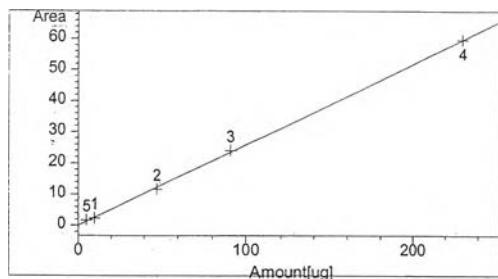
=====
 Peak Sum Table
 =====

No Entries in table

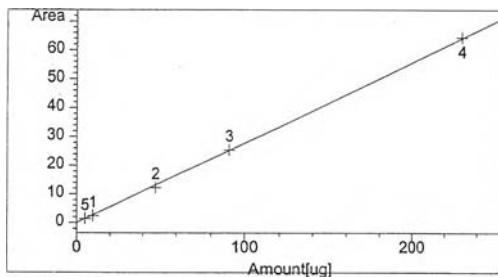
=====
 Calibration Curves
 =====



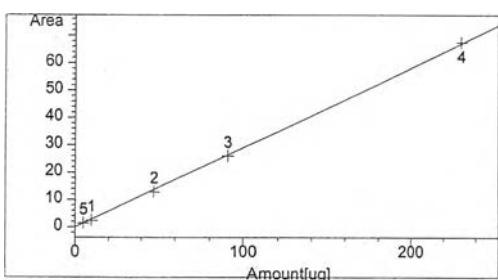
Method C:\HPCHEM\1\METHODS\CU2.M



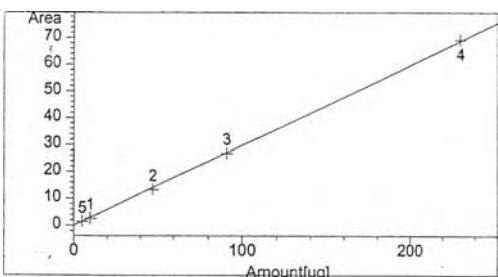
BENZENE at exp. RT: 8.945
 FID1 A,
 Correlation: 0.99989
 Residual Std. Dev.: 0.49171
 Formula: $y = mx$
 $m: 2.58886e-1$
 $x: \text{Amount [ug]}$
 $y: \text{Area}$



TOLUENE at exp. RT: 12.671
 FID1 A,
 Correlation: 0.99987
 Residual Std. Dev.: 0.55738
 Formula: $y = mx$
 $m: 2.77740e-1$
 $x: \text{Amount [ug]}$
 $y: \text{Area}$

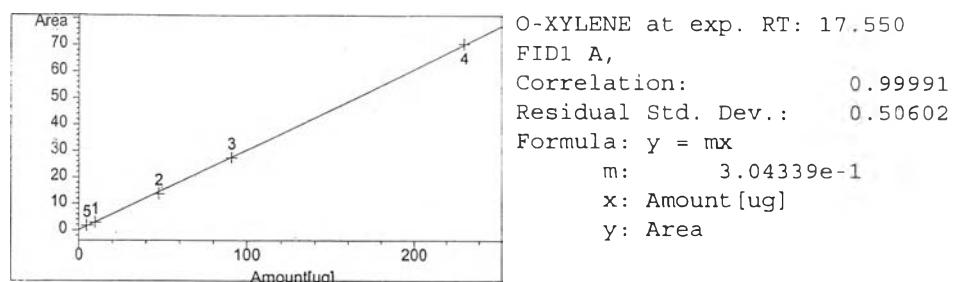


ETHYLBENZENE at exp. RT: 16.401
 FID1 A,
 Correlation: 0.99985
 Residual Std. Dev.: 0.63016
 Formula: $y = mx$
 $m: 2.90790e-1$
 $x: \text{Amount [ug]}$
 $y: \text{Area}$



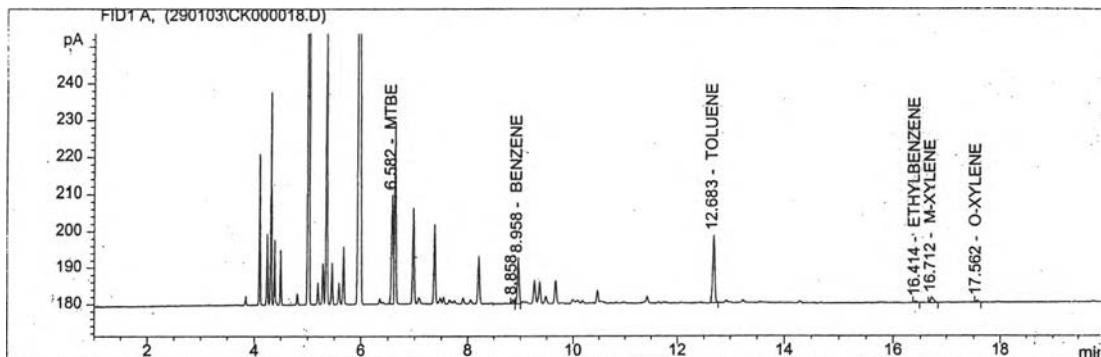
M-XYLENE at exp. RT: 16.696
 FID1 A,
 Correlation: 0.99989
 Residual Std. Dev.: 0.56488
 Formula: $y = mx$
 $m: 2.98511e-1$
 $x: \text{Amount [ug]}$
 $y: \text{Area}$

Method C:\HPCHEM\1\METHODS\CU2.M



Chromatogram of BTEX and MTBE

```
=====
Injection Date : 1/30/03 1:17:45 AM           Seq. Line : 18
Sample Name    : front7                      Vial   : 17
Acq. Operator   : Chanpen                   Inj    : 1
                                         Inj Volume : 1 µl
Acq. Method     : C:\HPCHEM\1\METHODS\CU2.M
Last changed    : 1/29/03 2:52:26 PM by Chanpen
Analysis Method : C:\HPCHEM\1\METHODS\CU2.M
Last changed    : 2/4/03 12:18:20 PM by somkiat
                                         (modified after loading)
```



```
Sorted By      : Signal
Calib. Data Modified : 2/4/03 11:15:14 AM
Multiplier     : 1.0000
Dilution       : 1.0000
```

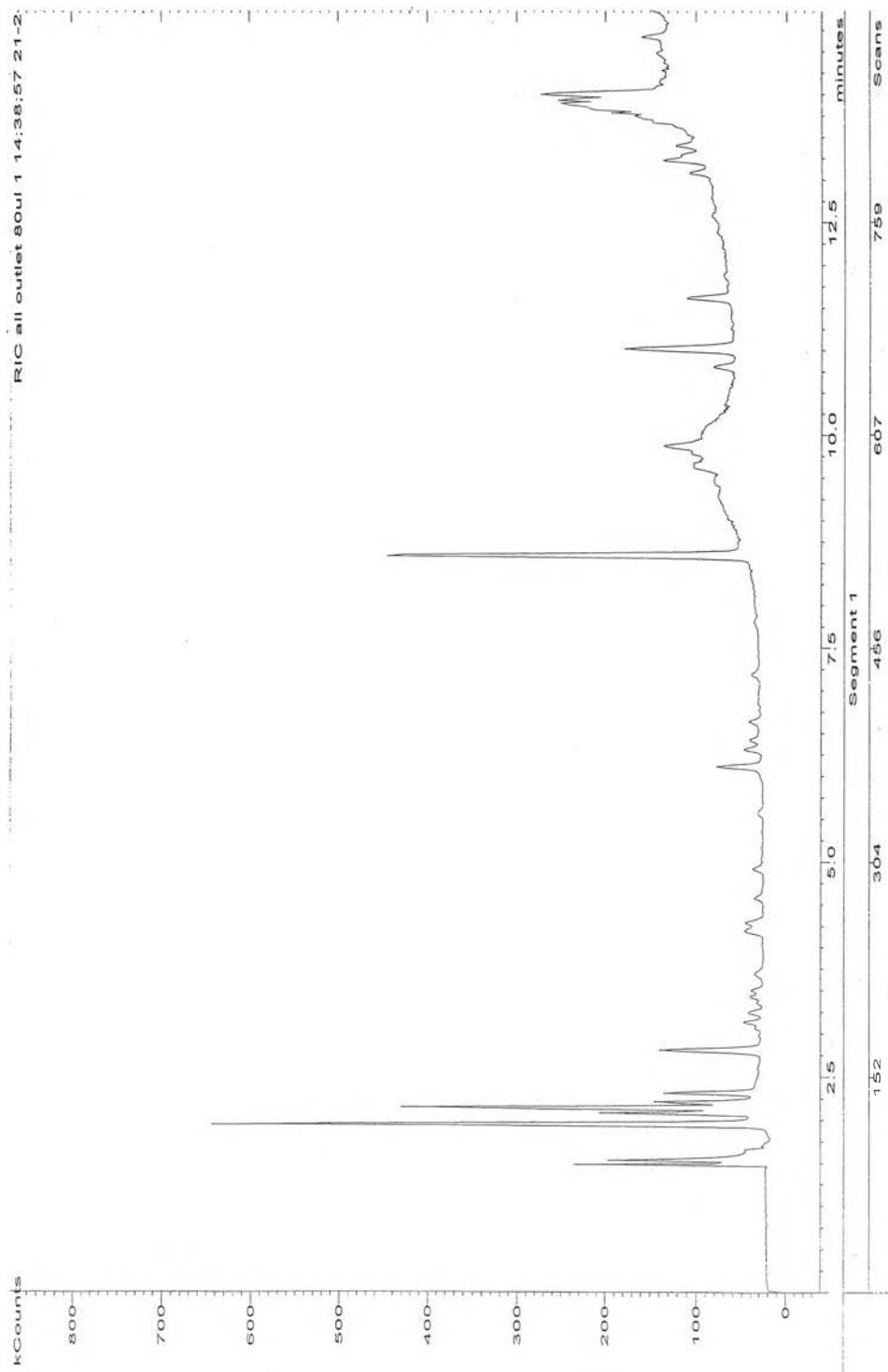
Signal 1: FID1 A,

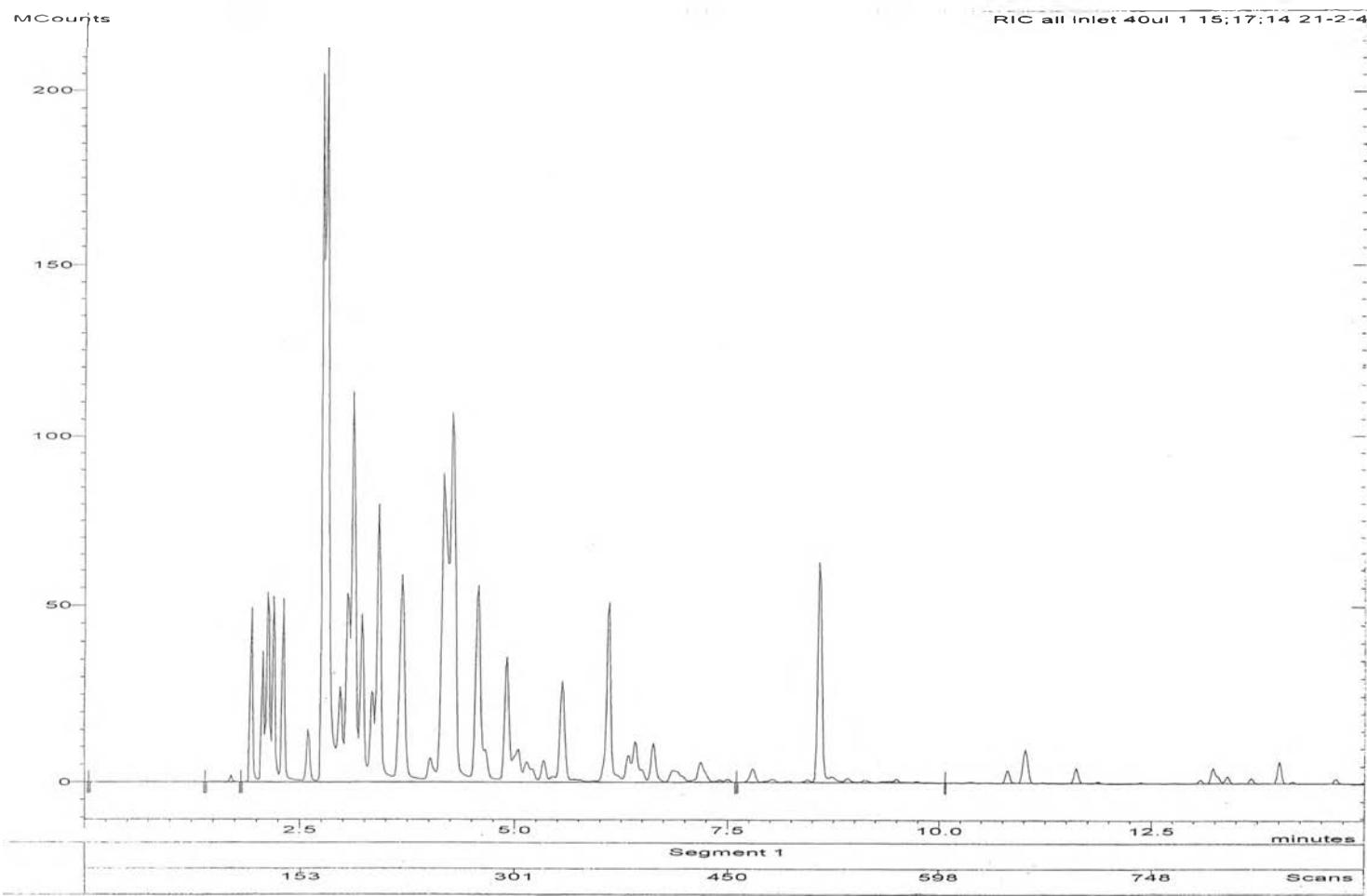
RetTime [min]	Type	Area [pA*s]	Amt/Area	Amount [ug]	Grp	Name
6.582	BV	63.10792	6.43377	406.02169		MTBE
8.958	VBA	28.40019	3.86270	109.70155		BENZENE
12.683	BB	47.19960	3.60049	169.94155		TOLUENE
16.414	BP	1.42584	3.43891	4.90332		ETHYLBENZENE
16.712	PB	6.41595	3.34997	21.49322		M-XYLENE
17.562	BP	1.67009	3.28581	5.48760		O-XYLENE

Totals : 717.54893

Results obtained with enhanced integrator!
1 Warnings or Errors :

APPENDIX B

APPENDIX B**Chromatogram of GC/MS**



Saturn Fit Search Hit List

Saturn Fit Search Hit List

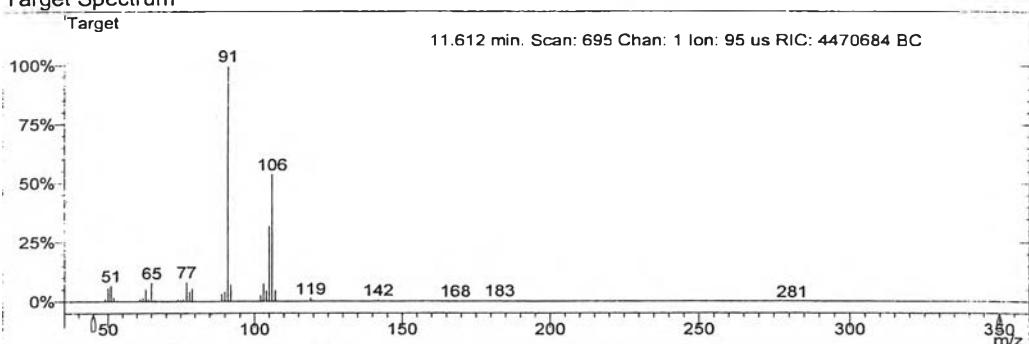
Saturn Fit Search Results

Hits Found: 25
 Pre-Search Hits Found: 350

Saturn Fit Search Parameters

Threshold: 700
 Target Ion Range: 45 - 350
 Library MW Range: 45 - 250
 Library Ion Range: 45 - 350 (Target Mass Range)
 Local Normalization: On
 Requested Pre-Search: 250
 Requested Final Search: 25
 Search 4 Libraries:
 A. c:\saturnws\satlib\nist98m.lbr
 B. c:\saturnws\satlib\nist98r.lbr
 C. c:\saturnws\satlib\libr_tr.lbr
 D. c:\saturnws\satlib\libr_tx.lbr

Target Spectrum



Spectrum from c:\...\cu oil\inlet 40ul 1 15;17;14 21-2-46.sms

Scan No: 695, Time: 11.612 minutes

No averaging. Background corrected.

Comment: 11.612 min. Scan: 695 Chan: 1 Ion: 95 us RIC: 4470684 BC

Pair Count: 67 MW: 0 Formula: None CAS No: None Acquired Range: 40 - 350

Purity	Fit	RFit	Entry #	MW.	Formula, CAS No., Name
1.	928	979	979	9522	B 106 C8H10, 95-47-6, Benzene, 1,2-dimethyl-
2.	944	977	977	9508	B 106 C8H10, 108-38-3, Benzene, 1,3-dimethyl-
3.	895	975	975	9512	B 106 C8H10, 2175-91-9, 1,3-Cyclopentadiene, 5-(
4.	927	975	975	9510	B 106 C8H10, 108-38-3, Benzene, 1,3-dimethyl-
5.	927	975	975	9517	B 106 C8H10, 106-42-3, p-Xylene
6.	938	971	971	9514	B 106 C8H10, 106-42-3, p-Xylene
7.	936	970	970	9515	B 106 C8H10, 106-42-3, p-Xylene
8.	936	969	969	37716	A 106 C8H10, 106-42-3, p-Xylene
9.	936	969	969	9509	B 106 C8H10, 108-38-3, Benzene, 1,3-dimethyl-
10.	936	969	969	9518	B 106 C8H10, 106-42-3, p-Xylene
11.	935	969	969	9523	B 106 C8H10, 95-47-6, Benzene, 1,2-dimethyl-
12.	935	968	968	9511	B 106 C8H10, 108-38-3, Benzene, 1,3-dimethyl-
13.	934	968	968	9516	B 106 C8H10, 106-42-3, p-Xylene
14.	934	967	967	9524	B 106 C8H10, 95-47-6, Benzene, 1,2-dimethyl-
15.	932	966	966	37714	A 106 C8H10, 108-38-3, Benzene, 1,3-dimethyl-
16.	932	965	965	37719	A 106 C8H10, 95-47-6, Benzene, 1,2-dimethyl-
17.	928	961	961	9500	B 106 C8H10, 95-47-6, Benzene, 1,2-dimethyl-
18.	889	956	956	37691	A 106 C8H10, 2175-91-9, 1,3-Cyclopentadiene, 5-(
19.	918	951	951	9513	B 106 C8H10, 2175-91-9, 1,3-Cyclopentadiene, 5-(
20.	451	944	944	36724	A 176 C9H8N2O2, 16844-42-1, Sydnone, 3-(phenylme
21.	899	941	941	36861	A 106 C8H10, 2809-71-4, 2,4-Octadiyne
22.	881	940	940	9354	B 106 C8H10, 2809-71-4, 2,4-Octadiyne
23.	409	939	939	36763	A 170 C7H7Br, 5376-03-4, Cycloheptatrienylium, b
24.	897	939	939	36408	A 106 C8H10, None, 1,6-Heptadien-3-yne, 5-methyl
25.	891	937	937	37696	A 106 C8H10, 100-41-4, Ethylbenzene

APPENDIX C

APPENDIX C

Daily calculation of VRU's ability for controlling gasoline vapor

Table C-1 Control Efficiency (CE%) of VRU at FPT

Test	MTBE	Benzene	Toluene	Ethybenzene	o-Xylene	m-Xylene	Total VOC	
21/1/03	test1	100.00	100.00	98.86	100.00	95.02	96.19	99.980
	test2	100.00	100.00	97.31	na	na	na	99.990
22/1/03	test1	100.00	99.91	99.44	100.00	100.00	100.00	99.991
	test2	100.00	100.00	99.74	100.00	100.00	100.00	99.996
23/1/03	test1	100.00	100.00	97.24	na	na	na	99.968
	test2	100.00	99.96	98.39	na	na	89.30	99.976
24/1/03	test1	100.00	100.00	100.00	na	na	100.00	99.984
	test2	100.00	100.00	98.71	na	na	na	99.996
25/1/03	test1	100.00	99.72	98.44	na	na	na	99.967
	test2	100.00	99.80	99.20	na	na	100.00	99.948
26/1/03	test1	100.00	100.00	100.00	100.00	100.00	100.00	99.989
	test2	100.00	100.00	100.00	na	100.00	100.00	99.999
27/1/03	test1	100.00	99.96	na	100.00	100.00	100.00	99.982
	test2	100.00	100.00	na	100.00	100.00	100.00	99.989
2/2/03	test1	100.00	error	100.00	na	na	100.00	99.978
	test2	100.00	error	100.00	na	100.00	100.00	99.974
Average		100.00	99.953	99.095	100.000	99.378	98.791	99.982
Average HAP ^a CE% 99.536								

error Analysis error

na Not found compound both inlet&outlet

^a Benzene, Toluene, Ethylbenzene, m-Xylene, o-Xylene and Methyl tert-Butyl Ether

Table C-2 Control Efficiency (CE%) of VRU at SHELL

Test	MTBE	Benzene	Toluene	Ethybenzene	o-Xylene	m-Xylene	Total VOC	
1/12/02	test1	99.739	98.751	na	100	na	98.013	99.981
	test2	99.892	95.792	na	na	100	na	99.994
2/12/02	test1	100	98.089	98.08	na	na	na	99.961
	test2	97.778	99.437	100	na	na	100	99.986
3/12/02	test1	100	98.756	na	na	na	na	99.961
	test2	100	92.957	na	na	na	na	99.983
5/12/02	test1	100	99.979	99.798	na	na	100	error
	test2	100	99.939	100	na	na	100	error
8/12/02	test1	99.856	95.525	na	na	na	na	99.983
	test2	100	99.495	na	na	na	na	99.977
12/12/02	test1	97.014	92.943	94.483	100	na	na	99.952
	test2	100	97.896	97.179	na	na	na	99.953
13/2/03	test1	99.927	98.198	98.542	100	na	95.778	99.902
	test2	99.98	99.189	99.173	100	100	97.352	99.96
14/2/03	test1	99.461	92.692	87.142	na	na	na	99.768
	test2	100	96.78	96.386	na	na	88.443	error
16/2/03	test1	100	99.92	99.343	100	100	98.081	99.991
	test2	99.954	99.539	97.763	100	100	95.92	99.976
Average	99.6445	97.548722	97.324083		100	100	97.065222	99.9552
Average HAP ^a CE% 98.597088								

error Analysis error

na Not found compound both inlet&outlet

^a Benzene, Toluene, Ethylbenzene, m-Xylene, o-Xylene and Methyl tert-Butyl Ether

BIOGRAPHY

Miss Chanpen Kanjnaprapan was born on February 7, 1974 in Satul, Thailand. She was graduated from the Occupational Health and Safety Department, Faculty of Public Health, Mahidol University in 1995.

She had worked at Sahaviriya Steel Industry (Public) Company as safety officer for two and haft years. In 1998, she had worked for Nestle Food (Thailand) as safety and environment supervisor for three and haft years. After that she started as a graduate student in the International Postgraduate Programs in Environmental Management, Inter-Department of Environmental Management, Chulalongkorn University in May 2001 and completed the program in May 2003.

