

Jim,
These field Blanks had
no internal standard or
surrogate added.

Quantitation Report

(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAY16\11410F.D

Vial: 16

Acq On : 17 May 2002 4:01 am

Operator:

Sample : nyc; fb

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: Lscint.p

Quant Time: May 17 04:43:06 2002

Quant Results File: W042302.RES

Quant Method : C:\MSDCHEM\1...\W042302.M (RTE Integrator)

Title : VOC method 8260

Last Update : Wed May 01 09:30:08 2002

Response via : Initial Calibration

DataAcq Meth : W042302

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	17.69	114	849	5.00	ug/L	0.00
27) Chlorobenzene-d5	25.47	117	1521	5.00	ug/L	0.00
51) 1,4-Dichlorobenzene-d4	31.56	150	1504	5.00	ug/L	0.02
System Monitoring Compounds						
2) 1,2-Dichloroethane-d4	16.76	65	397	12.85	ug/L	0.00
Spiked Amount 5.000			Recovery	=	257.00%	
17) Toluene-d8	21.62	98	1825	8.01	ug/L	0.00
Spiked Amount 5.000			Recovery	=	160.20%	
54) 4-Bromofluorobenzene	28.51	95	826	4.81	ug/L	0.00
Spiked Amount 5.000			Recovery	=	96.20%	
Target Compounds						
4) Chloromethane	6.19	50	1834	49.80	ug/L	79
11) Acetone	9.39	43	130676	55850.78	ug/L	98
13) Methylene Chloride	11.04	49	9188	308.24	ug/L	98
58) 1,3,5-Trimethylbenzene	29.23	105	1118	3.48	ug/L	71

 (#) = qualifier out of range (m) = manual integration

11410F.D W050302.M Fri May 24 15:09:32 2002

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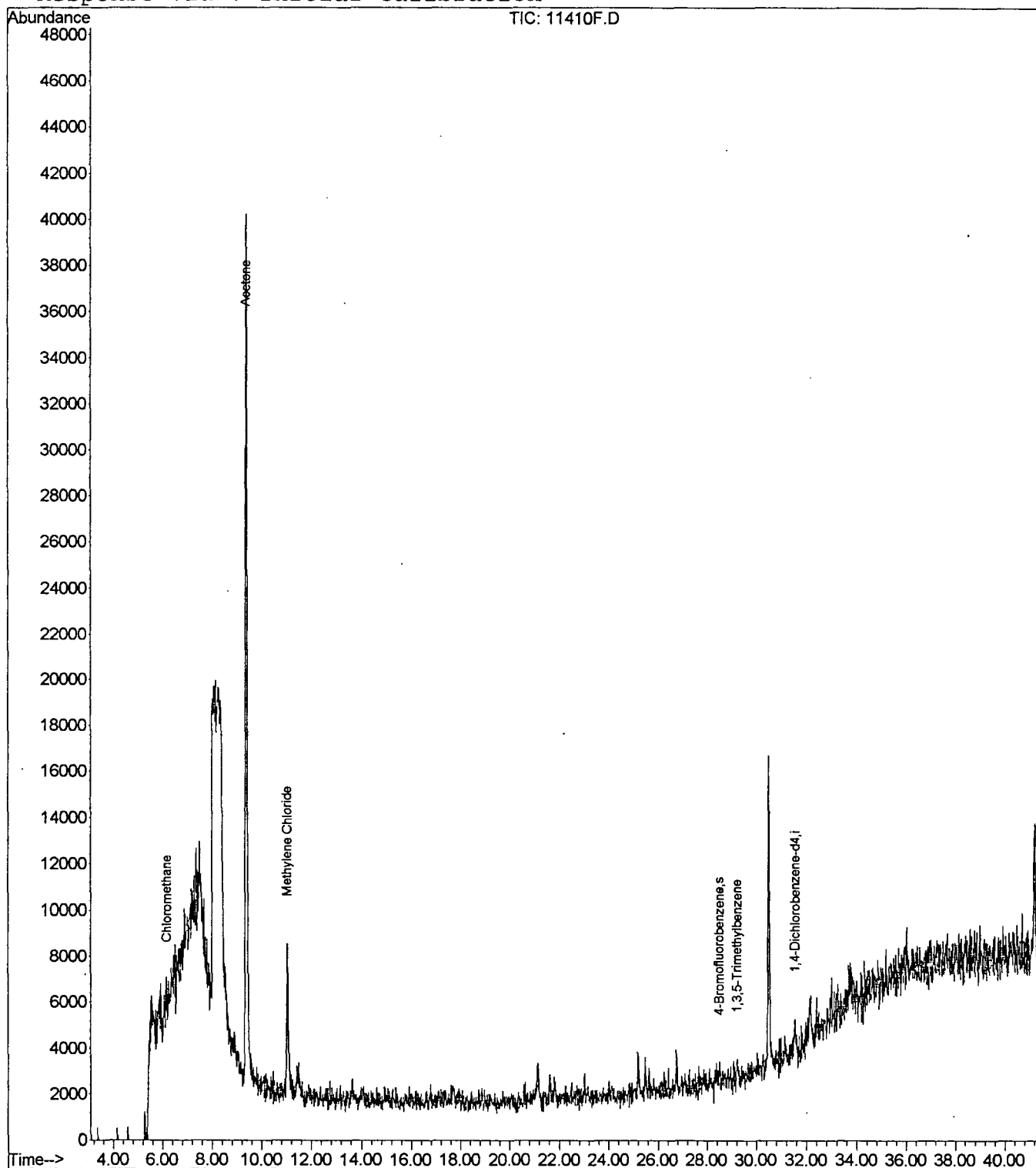
Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAY16\11410F.D
Acq On : 17 May 2002 4:01 am
Sample : nyc; fb
Misc :
MS Integration Params: Lscint.p
Quant Time: May 17 4:43 2002

Vial: 16
Operator:
Inst : Instrumen
Multiplr: 1.00

Quant Results File: W042302.RE

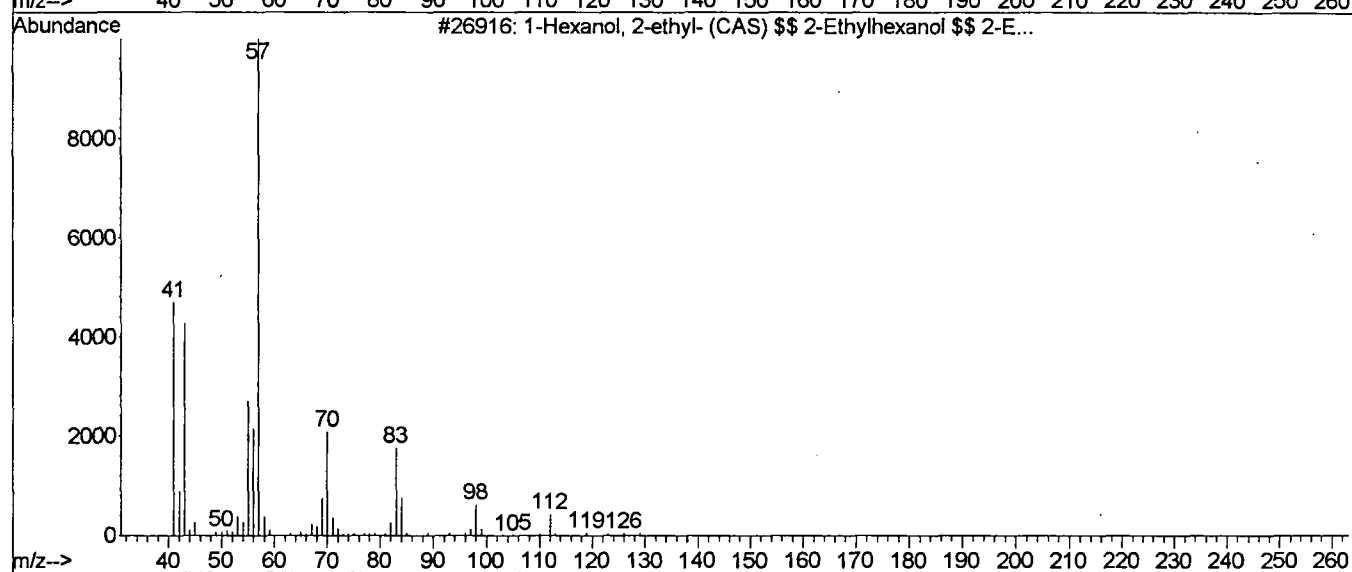
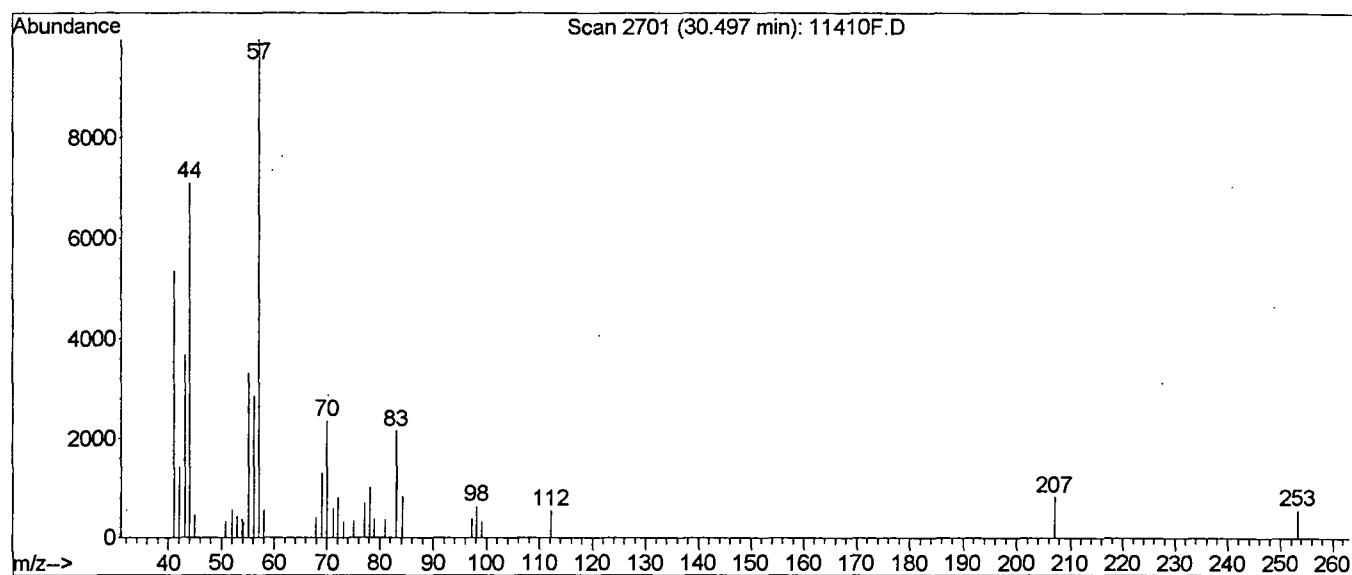
Method : C:\MSDCHEM\1\METHODS\METHODS\W050302.M (RTE Integrator)
Title : VOC method 8260
Last Update : Mon May 06 12:25:29 2002
Response via : Initial Calibration



Library Searched : C:\DATABASE\wiley7n.L

Quality : 47

ID : 1-Hexanol, 2-ethyl- (CAS) \$\$ 2-Ethylhexanol \$\$ 2-Ethyl-1-hexanol \$\$ Ethylhexanol \$\$ 2-Ethylhexan-1-ol
 \$\$ 2-Ethylhexyl alcohol \$\$ 2-Ethyl-hexanol-1 \$\$ Ethylhexyl alcohol \$\$ Octyl alcohol \$\$ 2-ETHYL-HEXAN-1-OL \$\$ 2-Ethylhexanol-1



Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAY16\11567F.D Vial: 18
 Acq On : 17 May 2002 5:34 am Operator:
 Sample : nyc; fb Inst : Instrumen
 Misc : Multiplr: 1.00
 MS Integration Params: Lscint.p
 Quant Time: May 17 06:16:03 2002 Quant Results File: W042302.RES

Quant Method : C:\MSDCHEM\1...\W042302.M (RTE Integrator)
 Title : VOC method 8260
 Last Update : Wed May 01 09:30:08 2002
 Response via : Initial Calibration
 DataAcq Meth : W042302

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	17.71	114	1277	5.00	ug/L	0.03
27) Chlorobenzene-d5	25.48	117	1254	5.00	ug/L	0.02
51) 1,4-Dichlorobenzene-d4	31.54	150	1187	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,2-Dichloroethane-d4	16.74	65	133	2.86	ug/L	-0.02
Spiked Amount 5.000			Recovery	=	57.20%	
17) Toluene-d8	21.63	98	1551	4.52	ug/L	0.02
Spiked Amount 5.000			Recovery	=	90.40%	
54) 4-Bromofluorobenzene	28.52	95	571	4.22	ug/L	0.02
Spiked Amount 5.000			Recovery	=	84.40%	
Target Compounds						Qvalue
4) Chloromethane	6.18	50	1957	35.33	ug/L	86
11) Acetone	9.40	43	11055	3141.30	ug/L	99
13) Methylene Chloride	11.05	49	4891	109.09	ug/L	88
28) Benzene	16.65	78	637	3.28	ug/L	95
36) Toluene	21.80	91	1511	6.52	ug/L	76
58) 1,3,5-Trimethylbenzene	29.22	105	1771	6.99	ug/L	66

 (#) = qualifier out of range (m) = manual integration

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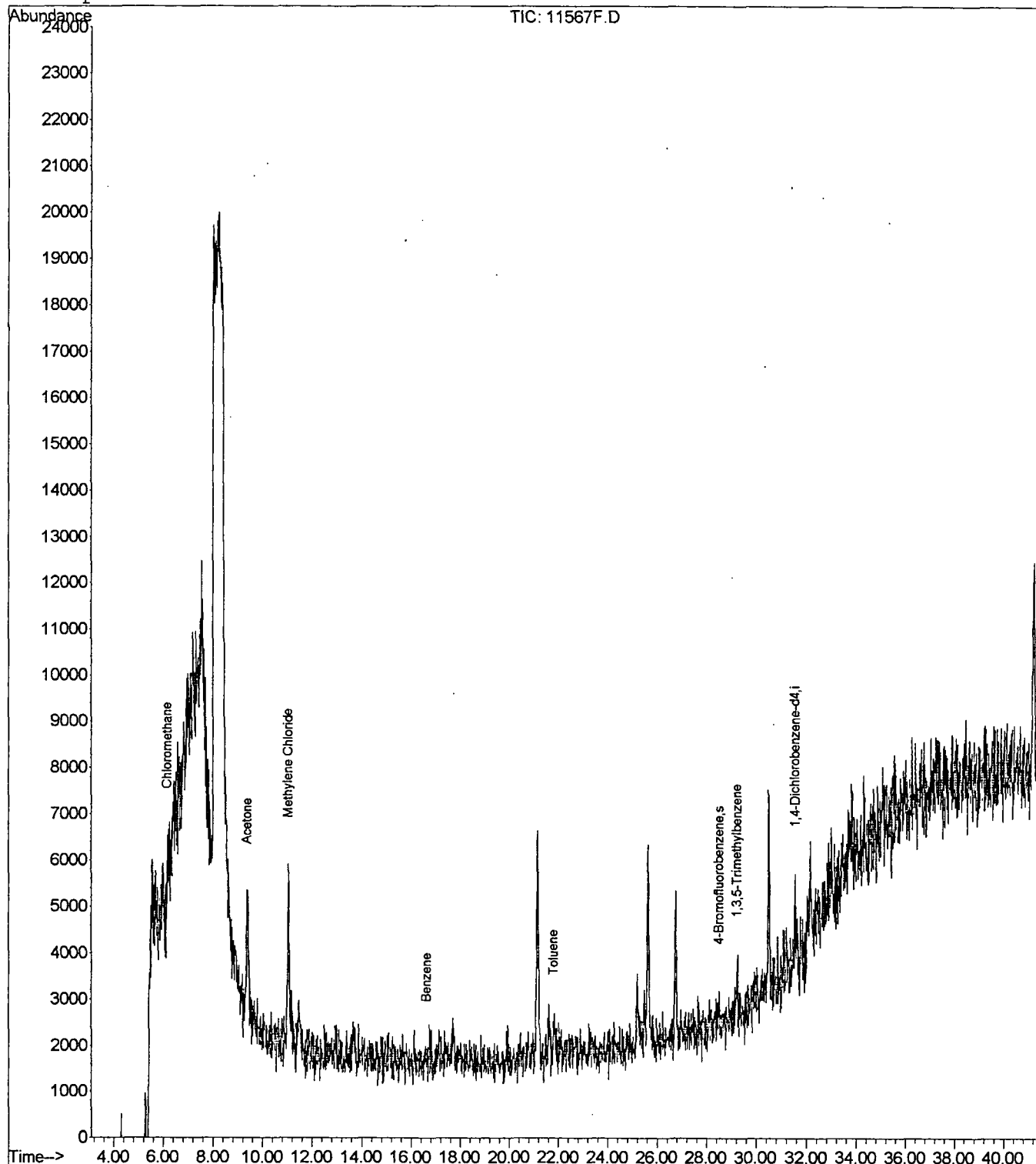
Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAY16\11567F.D
 Acq On : 17 May 2002 5:34 am
 Sample : nyc; fb
 Misc :
 MS Integration Params: Lscint.p
 Quant Time: May 17 6:16 2002

Vial: 18
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Results File: W042302.RE

Method : C:\MSDCHEM\1\METHODS\METHODS\W050302.M (RTE Integrator)
 Title : VOC method 8260
 Last Update : Mon May 06 12:25:29 2002
 Response via : Initial Calibration



Quantitation Report (Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAY16\11570F.D Vial: 19
 Acq On : 17 May 2002 6:21 am Operator:
 Sample : nyc; fb Inst : Instrumen
 Misc : Multiplr: 1.00
 MS Integration Params: Lscint.p
 Quant Time: May 17 07:02:35 2002 Quant Results File: W042302.RES

Quant Method : C:\MSDCHEM\1...\W042302.M (RTE Integrator)
 Title : VOC method 8260
 Last Update : Wed May 01 09:30:08 2002
 Response via : Initial Calibration
 DataAcq Meth : W042302

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	17.70	114	1272	5.00	ug/L	0.00
27) Chlorobenzene-d5	25.47	117	1285	5.00	ug/L	0.00
51) 1,4-Dichlorobenzene-d4	31.54	150	1253	5.00	ug/L	0.00
System Monitoring Compounds						
2) 1,2-Dichloroethane-d4	16.77	65	331	7.15	ug/L	0.00
Spiked Amount	5.000		Recovery	=	143.00%	
17) Toluene-d8	21.62	98	1685	4.93	ug/L	0.00
Spiked Amount	5.000		Recovery	=	98.60%	
54) 4-Bromofluorobenzene	28.50	95	311	2.18	ug/L	0.00
Spiked Amount	5.000		Recovery	=	43.60%	
Target Compounds						Qvalue
4) Chloromethane	5.67	50	1121	20.32	ug/L	97
7) Chloroethane	7.11	64	2291	66.31	ug/L	74
11) Acetone	9.40	43	17967	5125.42	ug/L	95
13) Methylene Chloride	11.05	49	9337	209.07	ug/L	97
28) Benzene	16.77	78	338	1.70	ug/L	97
36) Toluene	21.83	91	2012	8.47	ug/L	88
58) 1,3,5-Trimethylbenzene	29.24	105	2422	9.05	ug/L	85

 (#) = qualifier out of range (m) = manual integration

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Quantitation Report

(Not Reviewed)

Data File : C:\MSDCHEM\1\5973N\02MAY16\11570F.D

Vial: 19

Acq On : 17 May 2002 6:21 am

Operator:

Sample : nyc; fb

Inst : Instrumen

Misc :

Multiplr: 1.00

MS Integration Params: Lscint.p

Quant Time: May 17 7:02 2002

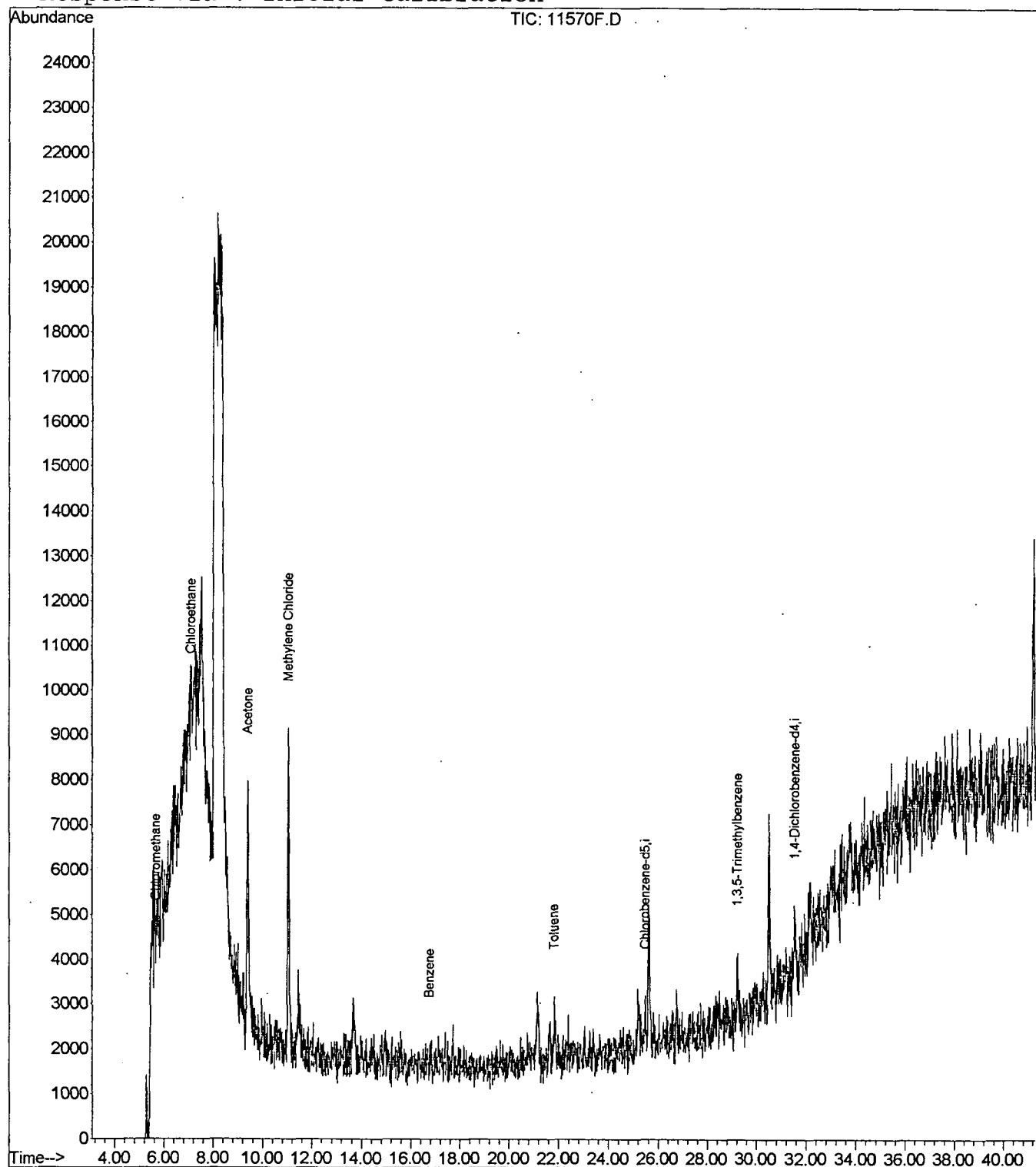
Quant Results File: W042302.RE

Method : C:\MSDCHEM\1\METHODS\METHODS\W050302.M (RTE Integrator)

Title : VOC method 8260

Last Update : Mon May 06 12:25:29 2002

Response via : Initial Calibration



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Library Searched : C:\DATABASE\wiley7n.L
 Quality : 74
 ID : Cyclotrisiloxane, hexamethyl-

