LAMPIRAN 1

DATA KARAKTERISASI KATALIS DENGAN BET

Quantachrome Corporation
Quantachrome Autosorb Automated Gas Sorption System Report
Autosorb for Windows© for AS-3 and AS-6 Version 1.23

Sample ID	Zeolit (2pk P-1)				
Description	B. Heru				
Comments					
Sample Weight	0.4379 g				
Adsorbate	NITROGEN	Outgas Temp	150.0 °C	Operator	Jajat
Cross-Sec Area	16.2 Å²/molecule	Outgas Time	2.0 hrs	Analysis Time	38.4 min
NonIdeality	6.580E-05	P/Po Toler	3	End of Run	06/26/2008 16:20
Molecular Wt	28.0134 g/mol	Equil Time	2	File Name	A8626024.RAW
Station #	4	Bath Temp.	77.40	PC SW Version	Pre-1.20

AREA-VOLUME-PORE SIZE SUMMARY

SURFACE AREA DATA

Multipoint BET	5.191E+01	m²/g
Langmuir Surface Area	8.051E+01	m²/g
t-Method External Surface Area	4.304E+01	m²/g
t-Method Micro Pore Surface Area	8.866E+00	m²/g
DR Method Micro Pore Area	7.137E+01	m²/g

PORE VOLUME DATA

t-Method Micro Pore Volume	4.538E-03	cc/g
DR Method Micro Pore Volume	2.536E-02	cc/g
HK Method Cumulative Pore Volume	2.207E-02	cc/g
SF Method Cumulative Pore Volume	2.249E-02	cc/g

PORE SIZE DATA

DR Method Micro Pore Width	9.687E+01	Å
DA Method Pore Diameter (Mode)	1.760E+01	Å
HK Method Pore Width (Mode)	1.402E+01	Å
SF Method Pore Diameter (Mode)	2.630E+01	Å

DATA REDUCTION PARAMETERS

Thermal Transpiration : OFF
Last Po Acquired 764.57 mm Hg
Additional Initialization Information Not Recorded.

BJH/DH Moving Average Size : 1

Interaction Constant (K) 2.9600 nm^3 x kJ/mol

Quantachrome Corporation
Quantachrome Autosorb Automated Gas Sorption System Report
Autosorb for Windows® for AS-3 and AS-6 Version 1.23

Sample ID	Zeolit + NH4 NO3				
Description	B. Heru				
Comments					
Sample Weight	0.3060 g				
Adsorbate	NITROGEN	Outgas Temp	150.0 °C	Operator	Jajat
Cross-Sec Area	16.2 Å²/molecule	Outgas Time	2.0 hrs	Analysis Time	59.0 min
NonIdeality	6.580E-05	P/Po Toler	3	End of Run	07/02/2008 14:52
Molecular Wt	28.0134 g/mol	Equil Time	2	File Name	A8702034.RAW
Station #	4	Bath Temp.	77.40	PC SW Version	Pre-1.20

AREA-VOLUME-PORE SIZE SUMMARY

SURFACE AREA DATA

Multipoint BET	1.016E+02	m²/g
Langmuir Surface Area	1.555E+02	m²/g
t-Method External Surface Area	6.318E+01	m²/g
t-Method Micro Pore Surface Area	3.844E+01	m²/g
DR Method Micro Pore Area	1.451E+02	m²/g

PORE VOLUME DATA

t-Method Micro Pore Volume	2.031E-02	cc/g
DR Method Micro Pore Volume	5.156E-02	cc/g
HK Method Cumulative Pore Volume	4.622E-02	cc/g
SF Method Cumulative Pore Volume	4.692E-02	cc/g

PORE SIZE DATA

DR Method Micro Pore Width	8.573E+01	Å
DA Method Pore Diameter (Mode)	1.740E+01	Å
HK Method Pore Width (Mode)	1.412E+01	Å
SF Method Pore Diameter (Mode)	2.639E+01	Å

DATA REDUCTION PARAMETERS

Thermal Transpiration : OFF
Last Po Acquired 765.66 mm Hg
Additional Initialization Information Not Recorded.

BJH/DH Moving Average Size : 1

Interaction Constant (K) 2.9600 nm^3 x kJ/mol

Quantachrome Corporation
Quantachrome Autosorb Automated Gas Sorption System Report
Autosorb for Windows® for AS-3 and AS-6 Version 1.23

Sample ID	Zeolit (HPA 5% O				
Description	B Heru				
Comments					
Sample Weight	0.4148 g				
Adsorbate	NITROGEN	Outgas Temp	250.0 °C	Operator	Jajat
Cross-Sec Area	16.2 Å²/molecule	Outgas Time	2.0 hrs	Analysis Time	64.6 min
NonIdeality	6.580E-05	P/Po Toler	3	End of Run	07/02/2008 10:32
Molecular Wt	28.0134 g/mol	Equil Time	2	File Name	A8702015.RAW
Station #	5	Bath Temp.	77.40	PC SW Version	Pre-1.20

AREA-VOLUME-PORE SIZE SUMMARY

SURFACE AREA DATA

Multipoint BET	6.371E+01	m²/g
Langmuir Surface Area	9.792E+01	m²/g
t-Method External Surface Area	4.166E+01	m²/g
t-Method Micro Pore Surface Area	2.205E+01	m²/g
DR Method Micro Pore Area	9.067E+01	m²/g

PORE VOLUME DATA

t-Method Micro Pore Volume	1.165E-02	cc/g
DR Method Micro Pore Volume	3.222E-02	cc/g
HK Method Cumulative Pore Volume	2.873E-02	cc/g
SF Method Cumulative Pore Volume	2.913E-02	cc/g

PORE SIZE DATA

DR Method Micro Pore Width	8.817E+01	Å
DA Method Pore Diameter (Mode)	1.740E+01	Å
HK Method Pore Width (Mode)	1.412E+01	Å
SF Method Pore Diameter (Mode)	2.657E+01	Å

DATA REDUCTION PARAMETERS

Thermal Transpiration : OFF
Last Po Acquired 736.16 mm Hg
Additional Initialization Information Not Recorded.

BJH/DH Moving Average Size : 1

Interaction Constant (K) 2.9600 nm^3 x kJ/mol

Quantachrome Corporation
Quantachrome Autosorb Automated Gas Sorption System Report
Autosorb for Windows© for AS-3 and AS-6 Version 1.23

Sample ID	Zeolit (HPAl0%)				
Description	B. Heru				
Comments					
Sample Weight	0.6077 g				
Adsorbate	Nitrogen	Outgas Temp	150.0 °C	Operator	Jajat
Cross-Sec Area	16.2 Å²/molecule	Outgas Time	2.0 hrs	Analysis Time	63.6 min
NonIdeality	6.580E-05	P/Po Toler	3	End of Run	06/26/2008 15:25
Molecular Wt	28.0134 g/mol	Equil Time	2	File Name	A8626016.RAW
Station #	6	Bath Temp.	77.40	PC SW Version	Pre-1.20

AREA-VOLUME-PORE SIZE SUMMARY

SURFACE AREA DATA

Multipoint BET	5.275E+01	m²/g
Langmuir Surface Area	8.259E+01	m²/g
t-Method External Surface Area	5.007E+01	m²/g
t-Method Micro Pore Surface Area	2.680E+00	m²/g
DR Method Micro Pore Area	7.242E+01	m²/g

PORE VOLUME DATA

t-Method Micro Pore Volume	1.211E-03	cc/g
DR Method Micro Pore Volume	2.574E-02	cc/g
HK Method Cumulative Pore Volume	2.170E-02	cc/g
SF Method Cumulative Pore Volume	2.218E-02	cc/g

PORE SIZE DATA

DR Method Micro Pore Width	1.082E+02	Å
DA Method Pore Diameter (Mode)	1.780E+01	Å
HK Method Pore Width (Mode)	1.398E+01	Å
SF Method Pore Diameter (Mode)	2.621E+01	Å

DATA REDUCTION PARAMETERS

Thermal Transpiration : ON
Effective Molecule Diameter (D) 3.5400 Å
Effective Cell Stem Inner Diameter (d) 4.0000 mm
Last Po Acquired 723.73 mm Hg
Additional Initialization Information Not Recorded.

BJH/DH Moving Average Size : 1

Interaction Constant (K) 2.9600 nm^3 x kJ/mol

Quantachrome Corporation
Quantachrome Autosorb Automated Gas Sorption System Report
Autosorb for Windows® for AS-3 and AS-6 Version 1.23

Sample ID	Zeolit HDA/ 20%				
Description	B. Heru				
Comments					
Sample Weight	0.3825 g				
Adsorbate	NITROGEN	Outgas Temp	150.0 °C	Operator	Jajat
Cross-Sec Area	16.2 Å²/molecule	Outgas Time	2.0 hrs	Analysis Time	57.2 min
NonIdeality	6.580E-05	P/Po Toler	3	End of Run	06/27/2008 12
Molecular Wt	28.0134 g/mol	Equil Time	2	File Name	A8627014.RAW
Station #	4	Bath Temp.	77.40	PC SW Version	Pre-1.20

AREA-VOLUME-PORE SIZE SUMMARY

SURFACE AREA DATA

Multipoint BET	2.428E+01	m²/g
Langmuir Surface Area	3.665E+01	m²/g
t-Method External Surface Area	1.161E+01	m²/g
t-Method Micro Pore Surface Area	1.267E+01	m²/g
DR Method Micro Pore Area	3.466E+01	m²/g

PORE VOLUME DATA

t-Method Micro Pore Volume	6.680E-03	cc/g
DR Method Micro Pore Volume	1.232E-02	cc/g
HK Method Cumulative Pore Volume	1.144E-02	cc/g
SF Method Cumulative Pore Volume	1.158E-02	cc/g

PORE SIZE DATA

DR Method Micro Pore Width	7.054E+01	Å
DA Method Pore Diameter (Mode)	1.640E+01	Å
HK Method Pore Width (Mode)	1.348E+01	Å
SF Method Pore Diameter (Mode)	2.528E+01	Å

DATA REDUCTION PARAMETERS

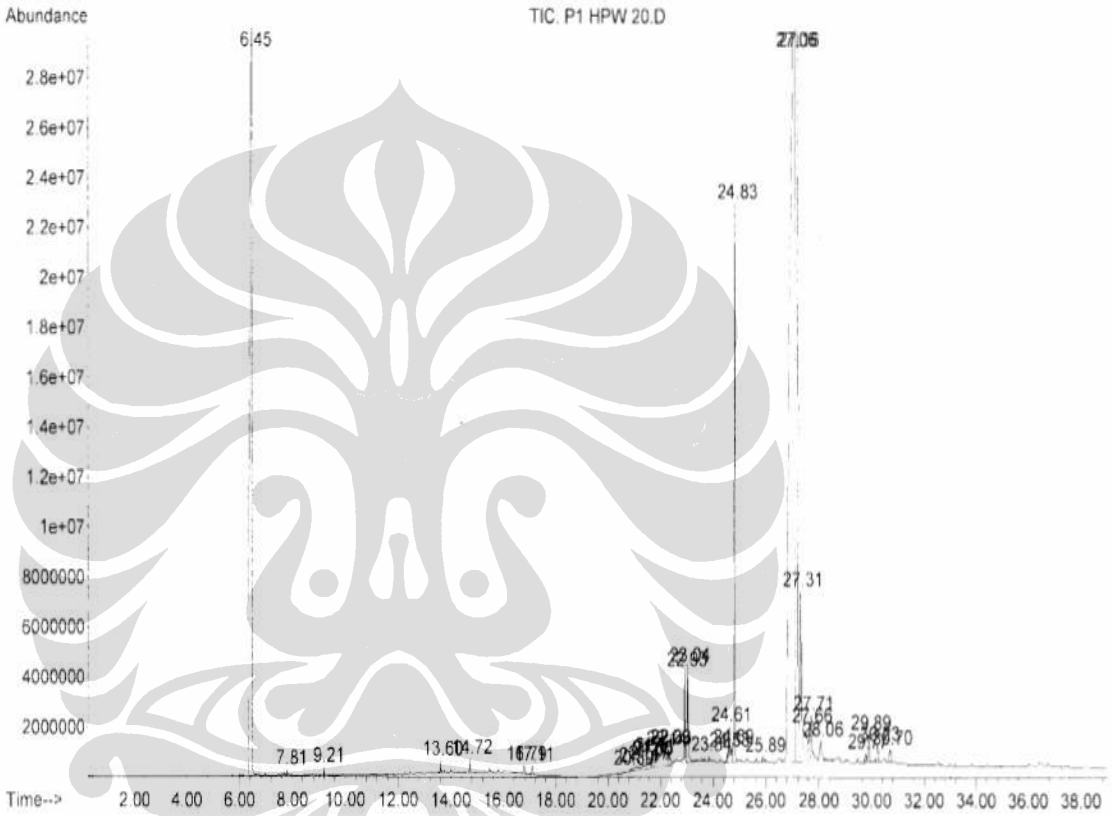
Thermal Transpiration : OFF
Last Po Acquired 763.80 mm Hg
Additional Initialization Information Not Recorded.

BJH/DH Moving Average Size : 1

Interaction Constant (K) 2.9600 nm^3 x kJ/mol

LAMPIRAN 2
DATA ANALISIS GC-MS

```
File       :C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\P1 HPW 20
...          .D
Operator   : PANJI DERMAWAN
Instrument :    Instrument #1
Acquired   :  2 Jul 2008   9:44     using AcqMethod ASAM LEMAK.M
Sample Name: P1 HPW /20
Misc Info  :
```



Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
1	6.45	14.76	C:\Database\wiley7n.l			
			1-Octanol	26811	000111-87-5	90
			1-Octanol (CAS) $$ Octilin $$ Alfol 8 $$ Sipol L8 $$ n-Octanol $$ n-Octan-1-ol $$ Octyl alcohol $$ Octanol $$ n-Octyl alcohol $$ Heptyl carbinol $$ Caprylic alcohol $$ 1-Hydroxyoctane $$ Octan-1-ol $$ Alcohol c-8 $$ Capryl alcohol $$ n-Heptyl carbinol $$	26818	000111-87-5	90
			1-Octanol (CAS) $$ Octilin $$ Alfol 8 $$ Sipol L8 $$ n-Octanol $$ n-Octan-1-ol $$ Octyl alcohol $$ Octanol $$ n-Octyl alcohol $$ Heptyl carbinol $$ Caprylic alcohol $$ 1-Hydroxyoctane $$ Octan-1-ol $$ Alcohol c-8 $$ Capryl alcohol $$ n-Heptyl carbinol $$	26814	000111-87-5	90
2	7.81	0.04	C:\Database\wiley7n.l			
			3-Decanol	58240	001565-81-7	80
			3-Decanol (CAS) $$ 3-HYDROXY-DECANE $$ 1-Ethyl-1-octanol $$ 3-Hydroxydecane	58242	001565-81-7	74
			3-Decanol	58243	001565-81-7	56
3	9.21	0.06	C:\Database\wiley7n.l			
			Tridecane	93371	000629-50-5	97
			Tridecane (CAS) $$ n-Tridecane $$ Tridecane, n-	93366	000629-50-5	97
			Tridecane (CAS) $$ n-Tridecane $$ Tridecane, n-	93373	000629-50-5	95
4	13.60	0.08	C:\Database\wiley7n.l			
			Heptanoic acid, octyl ester	176813	005132-75-2	72
			Cyclooctane	14045	000292-64-8	70
			Heptanoic acid, octyl ester	176811	005132-75-2	64
5	14.71	0.09	C:\Database\wiley7n.l			
			Octanoic acid, octyl ester $$ n-Octyl caprylate $$ Octyl caprylate $$ Octyl octanoate	195491	002306-88-9	90
			Cyclooctane (CAS) $$ Octamethylene $$ cyclooctane, octamethylene	14044	000292-64-8	38
			Cyclohexane, 1,2-dimethyl-, cis-	14022	002207-01-4	38
6	16.79	0.12	C:\Database\wiley7n.l			
			(E)-1,1-Dimethyl-2-(3,3-dimthyl-2-butylidene)-cyclopropane $$ Cyclopropane, 1,1-dimethyl-2-(1,2,2-trimethylpropylidene)-, (E)- (CAS)	50468	085851-40-7	46
			1-Cyclohexyl-2-methyl-prop-2-en-1-one	50314	025183-82-8	41

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
			1-Undecene, 8-methyl-	71436	074630-40-3	30
7	17.11	0.13	C:\Database\wiley7n.l			
			Decanoic acid, octyl ester	231406	002306-92-5	81
			Cyclooctane (CAS) $$ Octamethylene $$ cyclooctane, octamethylene	14044	000292-64-8	42
			Decanoic acid, hexyl ester	195537	010448-26-7	41
8	20.89	-0.07	C:\Database\wiley7n.l			
			4,14-BIS(HYDROXYMETHYL)-[2.2]METACYCLOPHANE $$ Tricyclo[9.3.1.1(4,8)]hexadeca-1(15),4,6,8(16),11,13-hexaene-5,14-dimethanol (CAS)	211269	056752-68-2	60
			3,5-Dimethylbenzaldehyde thiocarbamoylhydrazone	125320	000000-00-0	47
			Phoxim $$ 3,5-Dioxa-6-aza-4-phosphaoct-6-ene-8-nitrile, 4-ethoxy-7-phenyl-, 4-sulfide $$ Glyoxylonitrile, phenyl-, oxime O,O-diethyl phosphorothioate $$ Bayer 5621 $$ Bayer 77488 $$ Bayer 9053 $$ Baythion $$ Benzeneacetonitrile, .alpha.-[(diethoxyphosphi	246723	014816-18-3	44
9	21.11	0.18	C:\Database\wiley7n.l			
			4,14-BIS(HYDROXYMETHYL)-[2.2]METACYCLOPHANE $$ Tricyclo[9.3.1.1(4,8)]hexadeca-1(15),4,6,8(16),11,13-hexaene-5,14-dimethanol (CAS)	211269	056752-68-2	62
			1-Nonadecene	208861	018435-45-5	56
			1-Heptadecene (CAS) $$ Hexahydroaplotaxene	171779	006765-39-5	53
10	21.47	0.27	C:\Database\wiley7n.l			
			Silicone grease, Siliconfett	392047	000000-00-0	46
			3'-hydroxy-5'-propylphenyl 2,4-dihydroxy-6-pentylbenzoate	304792	000000-00-0	43
			2-Ethylacridine	125906	000000-00-0	42
11	21.60	0.17	C:\Database\wiley7n.l			
			13-Oxabicyclo[9.3.1]pentadecane, 1 5-chloro-	179051	000000-00-0	45
			2',4'-DIMETHYLOXANILIC ACID N'-VERATRYLIDENEHYDRAZIDE	302366	000000-00-0	44
			Hexahydropyridine, 1-methyl-4-[4,5-dihydroxyphenyl]-	125674	094427-47-1	43
12	21.65	0.09	C:\Database\wiley7n.l			
			9-Octadecenoic acid, (E)-	228774	000112-79-8	45
			HEPTADECENE-(8)-CARBONIC ACID-(1)	228686	000000-00-0	45
			Dodecahydropyrido[1,2-b]isoquinolin-6-one	125788	000000-00-0	43

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
13	21.71	0.16	C:\Database\wiley7n.l			
			Pentadecane, 2,6,10,14-tetramethyl - (CAS) $$ Pristane $$ PRISTANE (FIELD ION) $$ 2,6,10,14-Tetramethylpentadecane $$ Pristan $$ Norphytan $$ Norphytane $$ Bute hydrocarbon $$ 2,6,10,14-TETRAMETHYL PENTADECANE	211501	001921-70-6	41
			6-Ethyl-4-oxadecane-1,2-dithiol	167044	000000-00-0	38
			3-Methylheneicosane $$ Heneicosane, 3-methyl-	261321	006418-47-9	38
14	22.14	1.15	C:\Database\wiley7n.l			
			(2,3-Diphenylcyclopropyl)methyl phenyl sulfoxide, trans-	283396	000000-00-0	35
			Acetic acid, oxo((1-phenylethyl)amino)-, hydrazide $$ Semioxamazide, 5-(.alpha.-methylbenzyl)- $$ 5-(.alpha.-Methylbenzyl)semioxamazide $$ 5-(.alpha.-Phenylethyl)semioxamazide	125308	000093-95-8	15
			1,2-Dibenzoyl-4-cyano-1,2-dihydrocinnoline $$ 4-Cinnolinecarbonitrile, 1,2-dibenzoyl-1,2-dihydro- (CAS	310130	076051-75-7	12
15	22.29	0.15	C:\Database\wiley7n.l			
			9,12-Octadecadien-1-ol (CAS) $$ OCTADECA-9,12-DIEN-1-OL	208785	001577-52-2	96
			Z,E-3,13-Octadecadien-1-ol	208809	000000-00-0	91
			9,12-Octadecadienoic acid, methyl ester	243170	002462-85-3	84
16	22.33	0.21	C:\Database\wiley7n.l			
			1,9-Tetradecadiene	107091	112929-06-3	95
			Z-7-Hexadecen-1-ol acetate	228747	000000-00-0	94
			Z-9-Hexadecen-1-ol	174492	000000-00-0	93
17	22.93	0.63	C:\Database\wiley7n.l			
			Octane, 2-bromo- (CAS) $$ 2-Bromooctane $$ 2-Octyl bromide $$ sec-Octyl bromide $$ 1-Methylheptyl bromide	102199	000557-35-7	38
			Octane, 2-bromo-	102195	000557-35-7	38
			bis(2-Ethylhexyl) ether	177063	000000-00-0	37
18	23.04	0.76	C:\Database\wiley7n.l			
			2-phenylethynylquinoline $$ Quinoline, 2-(phenylethynyl)- (CAS) $$ 2-(Phenylethynyl)quinoline	158542	070437-00-2	80
			Cyclooctane (CAS) $$ Octamethylene $$ cyclooctane, octamethylene	14044	000292-64-8	38
			MYRISTIC ACID, OCTYL ESTER	290453	000000-00-0	38

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
19	23.84	0.04	C:\Database\wiley7n.l			
			Silicone grease, Siliconfett	392047	000000-00-0	53
			Octasiloxane, 1,1,3,3,5,5,7,7,9,9, 11,11,13,13,15,15-hexadecamethyl-	379834	019095-24-0	49
			1H-Indole, 1-methyl-2-phenyl- $$ I ndole, 1-methyl-2-phenyl- $$ 1-Met hyl-2-phenylindole $$ 2-Phenyl-N-m ethylindole $$ N-Methyl-2-phenylin dole	125864	003558-24-5	47
20	24.53	0.12	C:\Database\wiley7n.l			
			1-Octadecene	190564	000112-88-9	59
			9-Octadecenoic acid, (E)- $$ trans -.delta.(sup 9)-Octadecenoic acid $$ trans-.delta.9-Octadecenoic aci d $$ trans-Octadec-9-enoic acid $$ trans-Oleic acid $$ trans-9-Octad ecenoic acid $$ Elaidic acid	228773	000112-79-8	58
			1-Docosene	259091	001599-67-3	56
21	24.60	0.37	C:\Database\wiley7n.l			
			Hexadecanoic acid, octadecyl ester (CAS) $$ Stearyl palmitate $$ Oct adecyl hexadecanoate $$ Octadecyl palmitate $$ Palmitic acid, octade cyl ester	370242	002598-99-4	90
			9-Hexadecenoic acid, octadecyl est er	369875	000000-00-0	81
			1-Heptadecene (CAS) $$ Hexahydroap lotaxene	171778	006765-39-5	78
22	24.69	0.14	C:\Database\wiley7n.l			
			13-Hexacosyne	308305	034291-68-4	64
			Z,Z-6,24-Tritriacontadien-2-one	362581	000000-00-0	50
			Bicyclo[4.1.0]heptane, 7-butyl- $$ Norcarane, 7-butyl-	50481	018645-10-8	50
23	24.83	7.19	C:\Database\wiley7n.l			
			2-Formyl-5-[4-(methoxycarbonyl)ben zoyl]pyrrole	196238	000000-00-0	50
			7-PHENYL-2-AZAFLUOREN-9-ONE $$ 9H-Indeno[2,1-c]pyridin-9-one, 7-phen yl- (CAS)	196578	062397-37-9	50
			Benz[a]acridine, 1,5-dimethyl-	196606	055030-47-2	47
24	25.90	0.07	C:\Database\wiley7n.l			
			Z-8-Methyl-9-tetradecenoic acid	174170	000000-00-0	44
			2-Ethylacridine	125906	000000-00-0	38
			2-Methyl-7-phenylindole	125901	001140-08-5	38
25	27.06	39.84	C:\Database\wiley7n.l			
			9-Octadecenoic acid (Z)-, 2-hydrox y-1-(hydroxymethyl)ethyl ester $$	303308	003443-84-3	62

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
			Olein, 2-mono- $$ .beta.-Monoolein $$ Glycerol 2-monooleate $$ 2-Monoolein $$ 2-Monooleoylglycerol $$ 2-Oleoyl glycerol ether $$ 2-Oleoylglycerol			
			9-Octadecenoic acid, (E)-	228772	000112-79-8	58
			Oleic acid, 3-hydroxypropyl ester	290335	000821-17-0	53
26	27.16	27.63	C:\Database\wiley7n.l			
			Decyl oleate $$ 9-Octadecenoic acid (Z)-, decyl ester	343892	003687-46-5	93
			8-Octadecenoic acid, methyl ester, (E)- $$ Methyl trans-8-octadecenoate	245464	026528-50-7	87
			Octadec-9-enoic acid	228692	000000-00-0	83
27	27.31	2.81	C:\Database\wiley7n.l			
			Octadecanoic acid, octyl ester $$ Stearic acid, octyl ester $$ Octyl octadecanoate $$ Octyl stearate	330276	000109-36-4	49
			Octadecanoic acid, octyl ester (CAS) $$ Octyl stearate $$ Octyl octadecanoate $$ Stearic acid, octyl ester $$ OCTADECANSAEURE, OCTYLESTE	330277	000109-36-4	47
			3,3':5',3''-bis(dimethylene)-2,2':6',2''-terpyridine $$ 5,6,8,9-tetrahydroquino[8,7-b][1,10]phenanthroline	232391	096413-21-7	43
28	27.66	0.48	C:\Database\wiley7n.l			
			7-Pentadecyne	128099	022089-89-0	96
			9,12-Octadecadien-1-ol (CAS) $$ OCTADECA-9,12-DIEN-1-OL	208785	001577-52-2	95
			Cyclohexene, 4-(4-ethylcyclohexyl)-1-pentyl-	203503	000000-00-0	91
29	27.72	0.79	C:\Database\wiley7n.l			
			9-Octadecenoic acid (Z)- (CAS) $$ Oleic acid $$ Red oil $$ Oelsauere $$ Oleine 7503 $$ Pamolyn 100 $$ Emersol 211 $$ Vopcolene 27 $$ cis-Oleic acid $$ Wecoline OO $$ Z-9-Octadecenoic acid $$ cis-9-Octadecenoic acid $$ .delta.9-cis-Oleic acid $$ 9-Octadece	228698	000112-80-1	90
			9-Octadecenoic acid, (E)-	228774	000112-79-8	87
			HEPTADECENE-(8)-CARBONIC ACID-(1)	228686	000000-00-0	87
30	28.06	0.51	C:\Database\wiley7n.l			
			9,12-Octadecadienoic acid (Z,Z)- (CAS) $$ Linoleic acid $$ Linoleic $$ Unifac 6550 $$ Linolic acid $$ Telfairic acid $$ Grape seed oil $	226101	000060-33-3	94

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
			$ Polylin No. 515 $$ cis,cis-Linol eic acid $$ 9,12-Octadecadienoic acid $$ cis-9,cis-12-Octadecadieno ic acid $$ 9,12-O			
			Z-4-Nonadecen-1-ol acetate	275569	000000-00-0	80
			Z,Z-10,12-Hexadecadien-1-ol acetat	226137	000000-00-0	78
31	29.77	0.14	C:\Database\wiley7n.l			
			1H-Indole, 5-methyl-2-phenyl- (CAS) $$ 5-Methyl-2-phenylindole $$ 2-Phenyl-5-methylindole $$ Indole, 5 -methyl-2-phenyl-	125868	013228-36-9	46
			1H-Indole, 5-methyl-2-phenyl- $$ I ndole, 5-methyl-2-phenyl- $$ 2-Phe nyl-5-methylindole $$ 5-Methyl-2-p henylindole	125869	013228-36-9	46
			1,1,1,3,5,5,5-Heptamethyltrisiloxa ne $$ Bis(trimethylsiloxy)methylsi lane $$ Hydromethylsiloxane $$ Tri siloxane, 1,1,1,3,5,5,5-heptamethy l-	146455	001873-88-7	42
32	29.89	0.44	C:\Database\wiley7n.l			
			2,6-Bis(methylthio)-4-(2-naphthyl) pyridine $$ Pyridine, 2,6-bis(meth ylthio)-4-(2-naphthalenyl)- (CAS)	246236	122913-46-6	50
			3-(3,5-DIOXO-4-PHENYL-1,2,4-TRIAZO LIDINE-1-YL)-3,10-DIMETHYL-2-METHY LIDENE-TRICYCLO(5.2.1.1*4,10)UNDEC A-5,8-DIENE $$ 1,2,4-Triazolidine-3,5-dione, 4-phenyl-1-(3a,6,7,7a-t etrahydro-7a,9-dimethyl-8-methylen e-1,6-ethano -1H-inden-9-yl)- (CAS) $$ 1,6-Ethano-1	307107	064624-83-5	38
			3-Hydroxydiphenylamine	94131	000101-18-8	35
33	30.23	0.25	C:\Database\wiley7n.l			
			1,3-Dioxolane, 4-ethyl-5-octyl-2,2 -bis(trifluoromethyl)-, trans-	298033	038274-73-6	43
			1-Bromo-11-iodoundecane	305739	000000-00-0	38
			4-Trifluoroacetoxypentadecane	274929	000000-00-0	38
34	30.70	0.20	C:\Database\wiley7n.l			
			(+-)-dihydronaulafine $$ ((+-)-7,1 2,13,14-tetrahydrocyclopent[de]ind olo[2',3':3,4]prido[1,2-b][2,7]nap hthyridin-4[6H]-one $$ Cyclopent[d e]indolo[2',3':3,4]pyrido[1,2-b][2 ,7]naphthyridin-4(6H)-one, 7,12,13 ,14-tetrahydro- (CAS)	264162	122577-05-3	64
			Methamino,imide derivative of 4-O-carbomethoxylamellicolic anhydride $$ 1H-Benz[de]isoquinoline-1,3(2H	263938	041634-21-3	52

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P1 HPW 20.D
Acq On : 2 Jul 2008 9:44
Operator : PANJI DERMAWAN
Sample : P1 HPW /20
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.1 Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - JSW.E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
			)-dione, 6-methoxy-2,7-dimethyl-4, 9-bis(methylamino)- (CAS) 6,7,9-Trimethoxy-3-methylbenzo[g]i soquinoline-5,10-quinone $$ Benz[g]isoquinoline-5,10-dione, 6,7,9-tr imethoxy-3-methyl- (CAS) $$ Bostry coidin dimethyl ether	263906	077784-10-2	47

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File       :C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\P II HPW
...         2 20%.D
Operator   : PANJI DERMAWAN
Instrument :    Instrument #1
Acquired   :  2 Jul 2008  10:49     using AcqMethod ASAM LEMAK.M
Sample Name: P II HPW /2  20%
Misc Info  :
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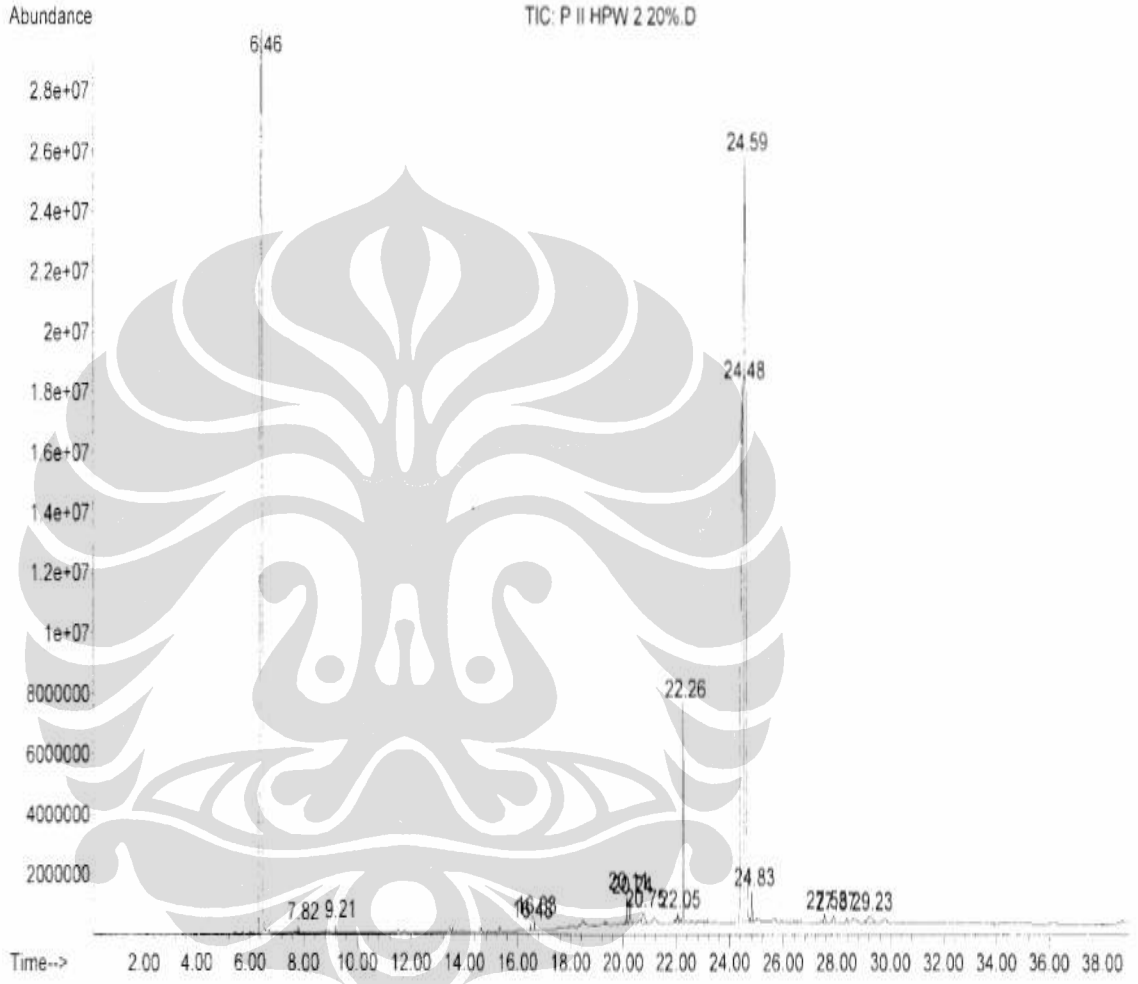

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P II HPW 2 20%.D
Acq On : 2 Jul 2008 10:49
Operator : PANJI DERMAWAN
Sample : P II HPW /2 20%
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - .E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
1	6.46	38.17	C:\Database\wiley7n.l			
			1-Octanol (CAS) $$ Octilin $$ Alfol 8 $$ Sipol L8 $$ n-Octanol $$ n-Octan-1-ol $$ Octyl alcohol $$ Octanol $$ n-Octyl alcohol $$ Heptyl carbinol $$ Caprylic alcohol $$ 1-Hydroxyoctane $$ Octan-1-ol $$ Alcohol c-8 $$ Capryl alcohol $$ n-Heptyl carbinol $$	26814	000111-87-5	90
			1-Octanol	26811	000111-87-5	90
			1-Octanol (CAS) $$ Octilin $$ Alfol 8 $$ Sipol L8 $$ n-Octanol $$ n-Octan-1-ol $$ Octyl alcohol $$ Octanol $$ n-Octyl alcohol $$ Heptyl carbinol $$ Caprylic alcohol $$ 1-Hydroxyoctane $$ Octan-1-ol $$ Alcohol c-8 $$ Capryl alcohol $$ n-Heptyl carbinol $$	26815	000111-87-5	90
	7.82	0.10	C:\Database\wiley7n.l			
			3-Decanol (CAS) $$ 3-HYDROXY-DECANE $$ 1-Ethyl-1-octanol $$ 3-Hydroxydecane	58242	001565-81-7	83
			3-Octanol, 6-ethyl-	58275	019781-27-2	83
			3-Decanol	58240	001565-81-7	80
	9.21	0.12	C:\Database\wiley7n.l			
			Tridecane	93371	000629-50-5	97
			Tridecane	93364	000629-50-5	93
			Tridecane	93367	000629-50-5	91
4	16.48	0.10	C:\Database\wiley7n.l			
			(E)-1,1-Dimethyl-2-(3,3-dimthyl-2-butylidene)-cyclopropane $$ Cyclopropane, 1,1-dimethyl-2-(1,2,2-trimethylpropylidene)-, (E)- (CAS)	50468	085851-40-7	50
			1-Cyclohexyl-2-methyl-prop-2-en-1-one	50314	025183-82-8	44
			Z-2-Octadecen-1-ol	211427	000000-00-0	41
5	16.64	0.13	C:\Database\wiley7n.l			
			Decanoic acid, octyl ester	231406	002306-92-5	55
			Decanoic acid, hexyl ester	195537	010448-26-7	46
			Cyclooctane	14045	000292-64-8	30
6	20.12	0.36	C:\Database\wiley7n.l			
			Octane, 2-bromo- (CAS) $$ 2-Bromooctane $$ 2-Bromoooctane $$ 2-Octyl bromide $$ sec-Octyl bromide $$ 1-Methylheptyl bromide	102199	000557-35-7	43
			Octane, 2-iodo- $$ sec-Octyl Iodide $$ 2-Iodooctane $$ 2-Octyl iodid	172825	000557-36-8	43
			Octane, 2-bromo-	102195	000557-35-7	43

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P II HPW 2 20%.D
Acq On : 2 Jul 2008 10:49
Operator : PANJI DERMAWAN
Sample : P II HPW /2 20%
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - .E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
7	20.24	0.46	C:\Database\wiley7n.l			
			2-phenylethynylquinoline $$ Quinoline, 2-(phenylethynyl)- (CAS) $$ 2-(Phenylethynyl)quinoline	158542	070437-00-2	64
			1-Nitrodibenzothiophene	159201	000000-00-0	49
			Octadecanoic acid, 11-methoxy-, methyl ester, (.+/-.)- $$ Octadecanoic acid, 11-methoxy-, methyl ester, (.+-.)-	279330	055334-49-1	25
8	20.76	0.50	C:\Database\wiley7n.l			
			1-Nonadecene	208860	018435-45-5	93
			Cyclohexane, 1,1'-(2-methyl-1,3-propanediyl)bis- $$ Propane, 1,3-dicyclohexyl-2-methyl- $$ 1,3-Dicyclohexyl-2-methylpropane	148908	002883-08-1	64
			Cyclohexane, 1,1'-(2-ethyl-1,3-propanediyl)bis-	168716	054833-34-0	64
9	22.05	0.20	C:\Database\wiley7n.l			
			Hexadecanoic acid, octadecyl ester (CAS) $$ Stearyl palmitate $$ Octadecyl hexadecanoate $$ Octadecyl palmitate $$ Palmitic acid, octadecyl ester	370242	002598-99-4	48
			Erucic acid $$ 13-Docosenoic acid, (Z)- $$ .delta.13-cis-Docosenoic acid $$ cis-13-Docosenoic acid	288671	000112-86-7	46
			9-Hexadecenoic acid, eicosyl ester, (Z)-	374599	022522-34-5	46
10	22.27	3.65	C:\Database\wiley7n.l			
			7-PHENYL-2-AZAFLUOREN-9-ONE $$ 9H-Indeno[2,1-c]pyridin-9-one, 7-phenyl- (CAS)	196578	062397-37-9	50
			Benz[c]acridine, 5,9-dimethyl- (CAS) $$ 5,9-DIMETHYLBENZO(A)ACRIDINE	196609	003518-03-4	47
			Benz[c]acridine, 5,9-dimethyl-	196610	003518-03-4	47
11	24.48	23.67	C:\Database\wiley7n.l			
			9,12-Octadecadienoic acid, methyl ester, (E,E)- (CAS) $$ Methyl linolelaidate $$ METHYL T9, T12 OCTADECADIENOATE $$ METHYL TRANS9, TRANS12-OCTADECADIENOATE $$ Linolelaidic acid, methyl ester $$ Methyl 9-trans-12-trans-octadecadienoate $$ Methyl trans,tran	243105	002566-97-4	86
			Isopropyl linoleate $$ 9,12-Octadecadienoic acid (Z,Z)-, 1-methylethyl ester	273683	022882-95-7	74
			1,3,12-Nonadecatriene	203500	000000-00-0	68

Library Search Report

Data Path : C:\MSDChem\1\DATA\MAHASISWA\S1 TK UI\PANJI DERMAWAN\
Data File : P II HPW 2 20%.D
Acq On : 2 Jul 2008 10:49
Operator : PANJI DERMAWAN
Sample : P II HPW /2 20%
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\wiley7n.l Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: Chemstation Integrator - .E

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
12	24.59	30.80	C:\Database\wiley7n.l			
			9-Octadecenoic acid (Z)-, hexadecyl ester $$ Oleic acid, hexadecyl ester $$ Cetyl oleate $$ Hexadecyl oleate $$ Palmityl oleate	369870	022393-86-8	93
			Oleic acid, eicosyl ester $$ 9-Octadecenoic acid (Z)-, eicosyl ester	378276	022393-88-0	90
			HEPTADECENE-(8)-CARBONIC ACID-(1)	228686	000000-00-0	87
13	24.83	0.61	C:\Database\wiley7n.l			
			Octadecanoic acid, octyl ester (CAS) $$ Octyl stearate $$ Octyl octadecanoate $$ Stearic acid, octyl ester $$ OCTADECANSAEURE, OCTYLESTE	330277	000109-36-4	50
			Octadecanoic acid, octyl ester	330275	000109-36-4	50
			3,3'-Diphenyl[2]staffane	234147	000000-00-0	47
14	27.53	0.24	C:\Database\wiley7n.l			
			5,6-Dimethoxy-2-(4'-methoxyphenyl)-3-methylindole	246379	000000-00-0	49
			2-(2'-Pyrazinyl)-phenanthro[9,10-d]oxazole	246402	000000-00-0	47
			Hexahydropyridine, 1-methyl-4-[4,5-dihydroxyphenyl]-	125674	094427-47-1	27
15	27.87	0.22	C:\Database\wiley7n.l			
			1,3-Dioxolane, 4-ethyl-5-octyl-2,2-bis(trifluoromethyl)-, trans-	298033	038274-73-6	38
			Cyclotrisiloxane, hexamethyl-	146396	000541-05-9	35
			2,4-Cyclohexadien-1-one, 3,5-bis(1,1-dimethylethyl)-4-hydroxy-	147861	054965-43-4	30
16	29.24	0.67	C:\Database\wiley7n.l			
			1-Nonadecene	208864	018435-45-5	70
			Hexadecane, 1-(ethenyloxy)- $$ Ether, hexadecyl vinyl $$ HEXADECAL VINYL ETHER $$ Cetyl vinyl ether $$ Vinyl cetyl ether $$ Hexadecyl vinyl ether	211413	000822-28-6	64
			1-Docosene	259091	001599-67-3	60

ASAM LEMAK.M Wed Jul 02 11:39:53 2008 Page: 1