

Method 508 mod

Date Extracted 2/5/02

Page _____

ref

Lab #	Description	Sample vol/wt	Drip rate ml/min	Surr add	Spk add	IS add	Final vol.	T.V. pos.	pH	Count
LB		500mL		1mL			1mL		7	
Ref					0.5mL					
2713										
2714										
2715										
2716										
2717										
2718										
2719										
2720										
2721										
Reviewed by, <i>[Signature]</i> 2/13/02										
Remarks All samples had a yellow tint after concentration BATCH# \$ GCMC- 17199										
drip rate must be between 5-15 ml/min for liq-liq extractors Extraction type - <u>sep fun</u> Solv.lot# <u>V41470</u> NaCl lot# <u>V39618</u> Na2SO4 lot# <u>V391100</u>										
STANDARD	INT STD LOG#									
CALIBRATION										
REF SPIKE	91759									
MDL/MDRF										
Q. CHECK										
INT. STD.										
SURROGATE	020102									
LPC										
ENDRIN/DDT										

Analyst LF/EL
 revision 4/12/01

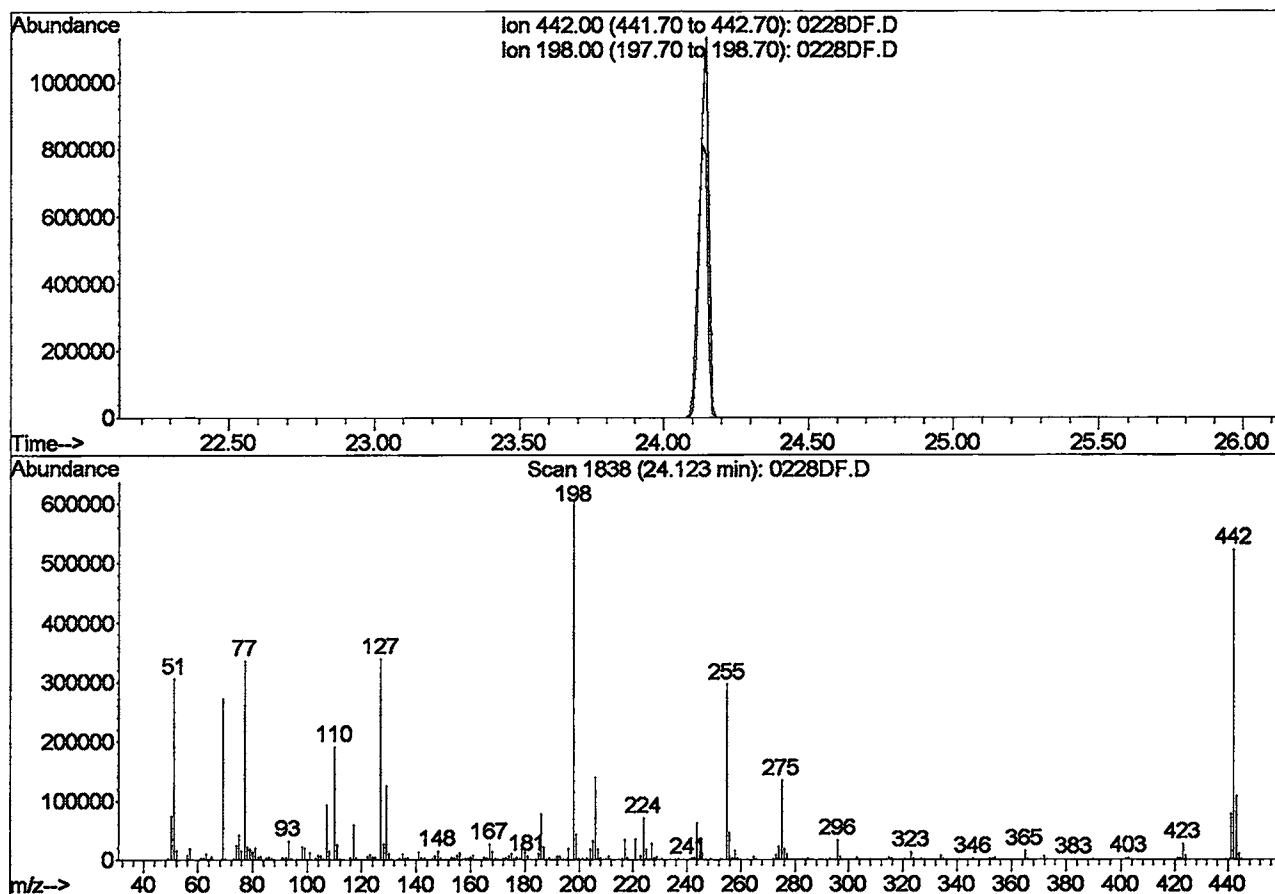
Verified by _____

Reviewed by *[Signature]*
 Date 3/5/02

DFTPP

Data File : C:\MSDCHEM\1\5973N\02FEB28\0228DF.D
 Acq On : 28 Feb 2002 8:42 am
 Sample : dftpp
 Misc : February 28, 2002
 MS Integration Params: SPLIT.P
 Method : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
 Title : 508 scan method

Vial: 1
 Operator:
 Inst : Instrume
 Multiplr: 1.00



Spectrum Information: Scan 1838

Target Mass	Rel. to Mass	Lower Limit%	Upper Limit%	Rel. Abn%	Raw Abn	Result Pass/Fail
51	198	30	60	50.4	305920	PASS
68	69	0.00	2	0.0	0	PASS
69	198	0.00	100	45.0	273152	PASS
70	69	0.00	2	0.4	1087	PASS
127	198	40	60	55.9	339584	PASS
197	198	0.00	1	0.0	0	PASS
198	198	100	100	100.0	607104	PASS
199	198	5	9	7.2	43752	PASS
275	198	10	30	22.2	134912	PASS
365	198	1	99	2.6	15935	PASS
441	443	1	99	71.3	77520	PASS
442	198	40	110	86.1	523008	PASS
443	442	17	23	20.8	108656	PASS

User: Baglioni, Walter
 Westchester Dept. of Labs & Research
 Date: 03/04/2002 Time: 12:15:59

Sample: AE02713
 Analysis: \$C1GCMS CALI GCMS
 Units: ug
 Started: 2/28/02 09:33 Ended: 3/4/02 09:33 Analyst: WB
 Result source: a:\0228c40.rr
 Page: 1

	Component Name	Viol	Result	Qualifier	MDL
1	n-Nitrosodimethylamine		41		2.0
2	bis(2-Chloroethyl)ether		48		2.0
3	1,3-Dichlorobenzene		38		2.0
4	1,4-Dichlorobenzene		38		2.0
5	1,2-Dichlorobenzene		42		2.0
6	2,2'-oxybis(1-Chloropropane)		45		2.0
7	n-Nitrosodi-n-propylamine		41		2.0
8	Hexachloroethane		39		2.0
9	Nitrobenzene		37		2.0
10	Isophorone		37		2.0
11	bis(2-Chloroethoxy)methane		39		2.0
12	1,2,4-Trichlorobenzene		37		2.0
13	Naphthalene		36		2.0
14	Hexachlorobutadiene		38		2.0
15	2-Chloronaphthalene		37		2.0
16	Dimethylphthalate		37		2.0
17	2,6-Dinitrotoluene		35		2.0
18	Acenaphthylene		36		2.0
19	Acenaphthene		36		2.0
20	2,4-Dinitrotoluene		33		2.0
21	Diethylphthalate		36		2.0
22	4-Chlorophenylphenylether		37		2.0
23	Fluorene		35		2.0
24	Azobenzene/1,2-Diphenylhydra		38		2.0
25	n-Nitrosodiphenylamine		30		2.0
26	4-Bromophenylphenylether		36		2.0
27	Hexachlorobenzene		38		2.0
28	Phenanthrene		34		2.0
29	Anthracene		35		2.0
30	Di-n-butylphthalate		36		2.0
31	Fluoranthene		35		2.0
32	Pyrene		36		2.0
33	Butylbenzylphthalate		38		2.0
34	Benzo(a)anthracene		37		2.0
35	bis(2-Ethylhexyl)phthalate		41		2.0
36	Chrysene		37		2.0
37	Di-n-octylphthalate		31		2.0
38	Benzo(b)fluoranthene		42		2.0
39	Benzo(k)fluoranthene		39		2.0
40	Benzo(a)pyrene		41		2.0
41	Indeno(1,2,3-cd)pyrene		43		2.0
42	Dibenz(a,h)anthracene		42		2.0
43	Benzo(g,h,i)perylene		39		2.0

Quantitation Report

(QT Reviewed)

Data File : C:\MSDCHEM\1\5973N\02FEB28\0228C40.D
 Acq Ok : 28 Feb 2002 9:33 am
 Sample : cal 40
 Misc : February 28, 2002
 MS Integration Params: SPLIT.P
 Quant Time: Mar 01 09:54:07 2002

Vial: 2
 Operator:
 Inst : Instrumen
 Multiplr: 1.00

Quant Results File: HP022102.RES

Quant Method : C:\MSDCHEM\1...\HP022102.M (RTE Integrator)

Title : 508 scan method

Last Update : Mon Feb 25 10:47:59 2002

Response via : Initial Calibration

DataAcq Meth : HP020702

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.72	152	1439437	20.00	ug/L	0.00
10) Naphthalene-d8	13.21	136	6420208	20.00	ug/L	0.00
17) Acenaphthene-d10	18.34	164	3292091	20.00	ug/L	0.00
29) Phenanthrene-d10	22.64	188	5508700	20.00	ug/L	0.00
38) Chrysene-d12	30.51	240	4401042	20.00	ug/L	-0.01
44) Perylene-d12	34.43	264	3409139	20.00	ug/L	-0.02

System Monitoring Compounds

37) 2,4-DDD	27.72	235	15349	0.21	ug/L	0.00
Spiked Amount	5.000		Recovery	=	4.20%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) n-Nitrosodimethylamine	3.98	74	2911078	41.26	ug/L	98
3) bis(2-Chloroethyl) ether	9.24	93	4506867	47.90	ug/L	99
4) 1,3-Dichlorobenzene	9.61	146	4072283	38.04	ug/L	98
5) 1,4-Dichlorobenzene	9.76	146	4046258	38.01	ug/L	99
6) 1,2-Dichlorobenzene	10.24	146	3948877	41.61	ug/L	100
7) 2,2'-oxybis (1-Chloropropa	10.68	45	6739512	45.10	ug/L	99
8) n-Nitroso-di-n-propylamine	11.06	70	2848789	40.58	ug/L	100
9) Hexachloroethane	11.03	117	1447243	39.31	ug/L	96
11) Nitrobenzene	11.37	77	4133778	36.62	ug/L	99
12) Isophorone	12.02	82	8014486	36.98	ug/L	99
13) bis(2-Chloroethoxy) methane	12.78	93	5488647	38.77	ug/L	99
14) 1,2,4-Trichlorobenzene	13.13	180	3098819	37.29	ug/L	98
15) Naphthalene	13.26	128	11333207	35.70	ug/L	100
16) Hexachlorobutadiene	13.85	225	1525764	37.62	ug/L	99
18) Hexachlorocyclopentadiene	15.96	237	1169580	44.96	ug/L	98
19) 2-Chloronaphthalene	16.66	162	6525046	37.08	ug/L	99
20) Dimethylphthalate	17.91	163	7328100	36.69	ug/L	99
21) 2,6-Dinitrotoluene	18.09	165	1591966	34.52	ug/L	96
22) Acenaphthylene	17.88	152	9576472	35.56	ug/L	100
23) Acenaphthene	18.44	153	6463166	36.38	ug/L	98
24) 2,4-Dinitrotoluene	19.20	165	2088175	32.79	ug/L	90
25) Diethylphthalate	20.01	149	6649830	35.77	ug/L	99
26) 4-Chlorophenyl-phenylether	20.05	204	3135330	37.30	ug/L	92
27) Fluorene	19.93	166	7473854	34.88	ug/L	100
28) Azobenzene/1,2-Diphenylhyd	20.49	77	9434664	38.49	ug/L	98
30) N-nitrosodiphenylamine	20.45	169	4159154	29.52	ug/L	99
31) 4-Bromophenyl-phenylether	21.45	248	2027758	36.47	ug/L	95
32) Hexachlorobenzene	21.80	284	2186721	37.74	ug/L	95
33) Phenanthrene	22.72	178	10604410	34.47	ug/L	100
34) Anthracene	22.86	178	10894235	35.14	ug/L	100

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

0228C40.D HP022102.M

Fri Mar 01 09:55:09 2002

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Data File : C:\MSDCHEM\1\5973N\02FEB28\0228C40.D
Acq On : 28 Feb 2002 9:33 am
Sample : cal 40
Misc : February 28, 2002
MS Integration Params: SPLIT.P
Quant Time: Mar 1 9:54 2002 Quant

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      Vial: 2
Operator:
Inst      : Instrumen
Multiplr: 1.00

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Quant Results File: HP022102.R

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Method      : C:\MSDCHEM\1\METHODS\METHODS\HP022102.M (RTE Integrator)
Title       : 508 scan method
Last Update : Mon Feb 25 10:47:59 2002
Response via : Initial Calibration
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