Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000850.D
Acq On : 04 Apr 2015 01:58
Operator : TP/IZ
Sample : MDL-02-W
Misc : MDL-WATER-0.2/0.4
ALS Vial : 39 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : MDL-02-W

Quant Time: Apr 04 04:21:02 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Apr 04 00:49:41 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	20076	5.00	ng		0.00
11) Naphthalene-d8	10.39	136	102621	5.00	ng		0.00
19) Acenaphthene-d10	14.26	164	52003	5.00	ng		0.00
26) Phenanthrene-d10	17.00	188	122237	5.00	ng		0.00
30) Chrysene-d12	21.21	240	96289	5.00	ng		0.00
34) Perylene-d12	23.39	264	88621	5.00	ng		-0.02
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	80796	14.61	ng		0.00
5) Phenol-d6	6.77	99	142004	13.33	ng		0.00
12) Nitrobenzene-d5	8.77	82	34939	7.30	ng		0.00
20) 2,4,6-Tribromophenol	15.77	330	20067	9.12	ng		0.00
23) 2-Fluorobiphenyl	12.87	172	160914	10.04	ng		0.00
32) Terphenyl-d14	19.65	244	132274	11.41	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	928	0.22	ng	#	66
3) n-Nitrosodimethylamine	3.46	42	366	0.14	ng	#	1
6) 2-Chlorophenol	7.17	128	1404	0.40	ng		85
7) Phenol	6.81	94	1520	0.18	ng		84
8) bis(2-Chloroethyl)ether	7.02	93	1105m	0.36	ng		
9) 2-Methylphenol	8.06	107	1052	0.31	ng	#	90
10) 3+4-Methylphenols	8.38	107	1088	0.26	ng		86
13) Nitrobenzene	8.80	77	1101	0.20	ng	#	69
14) 2-Nitrophenol	9.52	139	647	0.35	ng	#	44
15) 2,4-Dimethylphenol	9.58	122	832	0.14	ng		86
16) 2,4-Dichlorophenol	10.07	162	758	0.33	ng		84
17) Hexachlorobutadiene	10.72	225	721	0.18	ng	#	71
18) 4-Chloro-3-methylphenol	11.69	107	829	0.16	ng		93
21) 2,4,6-Trichlorophenol	12.70	196	350	0.21	ng	#	49
22) 2,4,5-Trichlorophenol	12.77	196	500	0.19	ng		84
25) 4-Nitrophenol	14.52	139	120	0.32	ng		90
27) 4,6-Dinitro-2-methylphenol	15.41	198	77	0.39	ng	#	86
28) Hexachlorobenzene	16.33	284	1031	0.20	ng	#	83
29) Pentachlorophenol	16.69	266	43	0.31	ng	#	67
31) Benzidine	19.29	184	1265	0.59	ng	#	72
33) 3,3'-Dichlorobenzidine	21.13	252	1062	0.31	ng	#	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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