Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040715\
Data File : BM000876.D
Acq On : 07 Apr 2015 12:51
Operator : TP/IZ
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC0.8

Quant Time: Apr 07 14:35:29 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 07 14:10:39 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	22690	5.00	ng	0.00
11) Naphthalene-d8	10.39	136	117869	5.00	ng	0.00
19) Acenaphthene-d10	14.26	164	63886	5.00	ng	0.00
26) Phenanthrene-d10	17.01	188	141268	5.00	ng	0.00
30) Chrysene-d12	21.20	240	126637	5.00	ng	0.00
34) Perylene-d12	23.38	264	113630	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	5929	0.93	ng	0.00
5) Phenol-d6	6.77	99	9644	1.07	ng	0.00
12) Nitrobenzene-d5	8.77	82	4470	0.81	ng	0.00
20) 2,4,6-Tribromophenol	15.75	330	1485	0.80	ng	0.00
23) 2-Fluorobiphenyl	12.87	172	15964	0.81	ng	0.00
32) Terphenyl-d14	19.65	244	12035	0.79	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.12	88	3063	0.86	ng	99
3) n-Nitrosodimethylamine	3.44	42	2735	0.94	ng	88
6) 2-Chlorophenol	7.17	128	7529	1.15	ng	98
7) Phenol	6.81	94	7210	0.77	ng	99
8) bis(2-Chloroethyl)ether	7.02	93	7614	1.12	ng	93
9) 2-Methylphenol	8.06	107	5532	0.83	ng	98
10) 3+4-Methylphenols	8.38	107	7291	1.01	ng	98
13) Nitrobenzene	8.80	77	4538	0.71	ng	92
14) 2-Nitrophenol	9.52	139	2508	0.75	ng	96
15) 2,4-Dimethylphenol	9.58	122	5216	0.78	ng	100
16) 2,4-Dichlorophenol	10.04	162	5078	0.85	ng	98
17) Hexachlorobutadiene	10.72	225	4339	0.93	ng	98
18) 4-Chloro-3-methylphenol	11.67	107	4902	0.83	ng	99
21) 2,4,6-Trichlorophenol	12.69	196	2596	0.79	ng	98
22) 2,4,5-Trichlorophenol	12.75	196	3432	0.84	ng	98
25) 4-Nitrophenol	14.51	139	790	0.66	ng	95
27) 4,6-Dinitro-2-methylphenol	15.40	198	1062	0.85	ng	89
28) Hexachlorobenzene	16.32	284	6287	0.79	ng	98
29) Pentachlorophenol	16.67	266	557	0.45	ng	96
31) Benzidine	19.27	184	4523	0.78	ng	97
33) 3,3'-Dichlorobenzidine	21.12	252	6693	0.77	ng	97

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

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