Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040915\
Data File : BM000930.D
Acq On : 09 Apr 2015 20:11
Operator : TP/IZ
Sample : SSTDICC0.8
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC0.8

Quant Time: Apr 10 06:07:44 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 10 05:53:53 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	40225	5.00	ng		0.00
8) Naphthalene-d8	10.38	136	152678	5.00	ng		0.00
17) Acenaphthene-d10	14.25	164	67150	5.00	ng		0.00
24) Phenanthrene-d10	16.99	188	154822	5.00	ng		0.00
30) Chrysene-d12	21.20	240	142899	5.00	ng		0.00
38) Perylene-d12	23.38	264	126396	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.21	112	6718	0.31	ng		0.00
5) Phenol-d6	6.78	99	8007	0.31	ng		0.00
9) Nitrobenzene-d5	8.76	82	5645	0.31	ng		0.00
18) 2,4,6-Tribromophenol	15.75	330	2364	0.32	ng		0.00
19) 2-Fluorobiphenyl	12.88	172	17715	0.32	ng		0.00
33) Terphenyl-d14	19.64	244	13007	0.30	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.12	88	3112	0.35	ng		97
3) n-Nitrosodimethylamine	3.43	42	3012	0.24	ng	#	81
6) Phenol	6.82	94	8169	0.32	ng		99
7) bis(2-Chloroethyl)ether	7.04	93	7447	0.34	ng		99
10) Nitrobenzene	8.80	77	6676	0.31	ng		100
11) 2,4-Dimethylphenol	9.56	122	6545	0.32	ng		99
12) 2,4-Dichlorophenol	10.03	162	5683	0.29	ng		95
13) Naphthalene	10.44	128	24922	0.33	ng		100
14) Hexachlorobutadiene	10.74	225	4556	0.31	ng		98
15) 2-Methylnaphthalene	12.06	142	16936	0.32	ng		99
16) 1-Methylnaphthalene	12.29	142	17109	0.33	ng		100
20) Acenaphthylene	13.96	152	28798	0.32	ng		99
21) Acenaphthene	14.32	154	16216	0.32	ng		99
22) 2,4-Dinitrophenol	14.40	184	446	0.32	ng		95
23) Fluorene	15.32	166	18035	0.31	ng		99
25) Hexachlorobenzene	16.32	284	7167	0.31	ng		99
26) Pentachlorophenol	16.67	266	1957	0.29	ng		98
27) Phenanthrene	17.04	178	33509	0.31	ng		99
28) Anthracene	17.14	178	29167	0.32	ng		99
29) Fluoranthene	19.07	202	34488	0.32	ng		99
31) Benzidine	19.26	184	5472	0.32	ng		99
32) Pyrene	19.43	202	36267	0.30	ng		100
34) Benzo(a)anthracene	21.18	228	31551	0.31	ng		100
35) 3,3'-Dichlorobenzidine	21.12	252	7681	0.31	ng		99
36) Chrysene	21.23	228	33214	0.31	ng		100
37) Indeno(1,2,3-cd)pyrene	25.56	276	33341	0.31	ng		100
39) Benzo(b)fluoranthene	22.72	252	30797	0.33	ng		99
40) Benzo(k)fluoranthene	22.76	252	31217	0.33	ng		99
41) Benzo(a)pyrene	23.28	252	27575	0.32	ng		99
42) Dibenzo(a,h)anthracene	25.57	278	25699	0.32	ng		99
43) Benzo(g,h,i)perylene	26.22	276	28538	0.32	ng		99

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

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