Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040915\
Data File : BM000931.D
Acq On : 09 Apr 2015 20:47
Operator : TP/IZ
Sample : SSTDICCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICCC001

Quant Time: Apr 10 06:08:23 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	39197	5.00	ng	0.00
8) Naphthalene-d8	10.38	136	149015	5.00	ng	0.00
17) Acenaphthene-d10	14.25	164	63813	5.00	ng	0.00
24) Phenanthrene-d10	16.99	188	147490	5.00	ng	0.00
30) Chrysene-d12	21.20	240	133532	5.00	ng	0.00
38) Perylene-d12	23.38	264	121474	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	8297	0.39	ng	0.00
5) Phenol-d6	6.78	99	9686	0.39	ng	0.00
9) Nitrobenzene-d5	8.76	82	6835	0.38	ng	0.00
18) 2,4,6-Tribromophenol	15.75	330	2812	0.39	ng	0.00
19) 2-Fluorobiphenyl	12.88	172	20859	0.39	ng	0.00
33) Terphenyl-d14	19.64	244	15897	0.40	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.12	88	3774	0.44	ng	100
3) n-Nitrosodimethylamine	3.43	42	3730	0.30	ng	100
6) Phenol	6.82	94	9931	0.40	ng	100
7) bis(2-Chloroethyl)ether	7.04	93	8886	0.41	ng	100
10) Nitrobenzene	8.80	77	8116	0.38	ng	100
11) 2,4-Dimethylphenol	9.56	122	7923	0.40	ng	100
12) 2,4-Dichlorophenol	10.03	162	6781	0.36	ng	100
13) Naphthalene	10.44	128	29321	0.39	ng	100
14) Hexachlorobutadiene	10.74	225	5314	0.37	ng	100
15) 2-Methylnaphthalene	12.06	142	20054	0.39	ng	100
16) 1-Methylnaphthalene	12.29	142	20141	0.40	ng	100
20) Acenaphthylene	13.96	152	34477	0.40	ng	100
21) Acenaphthene	14.32	154	18927	0.39	ng	100
22) 2,4-Dinitrophenol	14.40	184	523	0.40	ng	100
23) Fluorene	15.32	166	21470	0.39	ng	100
25) Hexachlorobenzene	16.32	284	8387	0.38	ng	100
26) Pentachlorophenol	16.67	266	2269	0.35	ng	100
27) Phenanthrene	17.04	178	40364	0.40	ng	100
28) Anthracene	17.14	178	33784	0.39	ng	100
29) Fluoranthene	19.07	202	41240	0.40	ng	100
31) Benzidine	19.26	184	6600	0.42	ng	100
32) Pyrene	19.43	202	43917	0.39	ng	100
34) Benzo(a)anthracene	21.18	228	37372	0.40	ng	100
35) 3,3'-Dichlorobenzidine	21.12	252	9420	0.41	ng	100
36) Chrysene	21.23	228	39758	0.40	ng	100
37) Indeno(1,2,3-cd)pyrene	25.56	276	39845	0.40	ng	100
39) Benzo(b)fluoranthene	22.72	252	36088	0.40	ng	100
40) Benzo(k)fluoranthene	22.76	252	37467	0.41	ng	100
41) Benzo(a)pyrene	23.28	252	32746	0.40	ng	100
42) Dibenzo(a,h)anthracene	25.57	278	30661	0.40	ng	100
43) Benzo(g,h,i)perylene	26.22	276	33980	0.40	ng	100

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

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