Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061315\
Data File : BM001682.D
Acq On : 12 Jun 2015 22:45
Operator : TP/IZ
Sample : SSTDICC0.5
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC0.5

Quant Time: Jun 13 00:28:34 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jun 13 00:26:17 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	14436	5.00	ng		0.00
6) Naphthalene-d8	10.46	136	52333	5.00	ng		0.00
12) Acenaphthene-d10	14.32	164	25549	5.00	ng		0.00
18) Phenanthrene-d10	17.06	188	79729	5.00	ng		0.00
25) Chrysene-d12	21.25	240	44224	5.00	ng		0.00
32) Perylene-d12	23.48	264	40517	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	1614	0.47	ng		0.00
5) Phenol-d6	6.83	99	2011	0.46	ng		0.00
7) Nitrobenzene-d5	8.83	82	1882	0.47	ng		0.00
13) 2,4,6-Tribromophenol	15.80	330	267	0.30	ng		0.00
14) 2-Fluorobiphenyl	12.94	172	4264	0.47	ng		0.00
27) Terphenyl-d14	19.70	244	3075	0.46	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.18	88	831	0.46	ng		97
3) n-Nitrosodimethylamine	3.49	42	1192	0.47	ng	#	99
8) Nitrobenzene	8.87	77	2000	0.46	ng		100
9) Naphthalene	10.51	128	5981	0.47	ng		98
10) Hexachlorobutadiene	10.78	225	1021	0.48	ng		99
11) 2-Methylnaphthalene	12.13	142	3806	0.46	ng		99
15) Acenaphthylene	14.03	152	5487	0.46	ng		100
16) Acenaphthene	14.39	154	3514	0.46	ng		99
17) Fluorene	15.36	166	5244	0.63	ng		100
19) 4-Bromophenyl-phenylether	16.24	248	693	0.10	ng		93
20) Hexachlorobenzene	16.38	284	1390	0.21	ng	#	100
21) Pentachlorophenol	16.72	266	509	0.27	ng		94
22) Phenanthrene	17.09	178	9352	0.37	ng		100
23) Anthracene	17.19	178	6609	0.23	ng		100
24) Fluoranthene	19.13	202	6626	0.19	ng		100
26) Pyrene	19.49	202	6867	0.45	ng		99
28) Benzo(a)anthracene	21.24	228	6450	0.46	ng		99
29) Chrysene	21.28	228	6228	0.47	ng		97
30) Bis(2-ethylhexyl)phthalate	21.16	149	3308	0.38	ng		100
31) Indeno(1,2,3-cd)pyrene	25.71	276	6288	0.57	ng		100
33) Benzo(b)fluoranthene	22.80	252	5788	0.44	ng		97
34) Benzo(k)fluoranthene	22.84	252	5579	0.43	ng		96
35) Benzo(a)pyrene	23.38	252	5053	0.46	ng		96
36) Dibenzo(a,h)anthracene	25.73	278	5058	0.53	ng		98
37) Benzo(g,h,i)perylene	26.40	276	5320	0.52	ng		98

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

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