Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061315\
Data File : BM001693.D
Acq On : 13 Jun 2015 05:24
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001EC

Quant Time: Jun 13 06:49:56 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 02:11:48 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	13525	5.00	ng		0.00
6) Naphthalene-d8	10.46	136	49519	5.00	ng		0.00
12) Acenaphthene-d10	14.32	164	24015	5.00	ng		0.00
18) Phenanthrene-d10	17.06	188	74054	5.00	ng		0.00
25) Chrysene-d12	21.25	240	42838	5.00	ng		0.00
32) Perylene-d12	23.48	264	39291	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	3279	1.01	ng		0.00
5) Phenol-d6	6.83	99	4136	1.01	ng		0.00
7) Nitrobenzene-d5	8.83	82	3855	0.99	ng		0.00
13) 2,4,6-Tribromophenol	15.80	330	599	1.07	ng		0.00
14) 2-Fluorobiphenyl	12.94	172	8591	1.01	ng		0.00
27) Terphenyl-d14	19.70	244	6252	0.98	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.18	88	1598	1.09	ng		99
3) n-Nitrosodimethylamine	3.49	42	2428	0.99	ng	#	98
8) Nitrobenzene	8.87	77	4141	0.98	ng		100
9) Naphthalene	10.51	128	11845	0.99	ng		99
10) Hexachlorobutadiene	10.78	225	1973	0.98	ng		98
11) 2-Methylnaphthalene	12.13	142	7586	0.99	ng		97
15) Acenaphthylene	14.03	152	11414	0.99	ng		100
16) Acenaphthene	14.39	154	6964	0.99	ng		98
17) Fluorene	15.36	166	9677	0.93	ng		99
19) 4-Bromophenyl-phenylether	16.24	248	1698	1.21	ng		79
20) Hexachlorobenzene	16.38	284	2397	0.90	ng	#	100
21) Pentachlorophenol	16.72	266	1115	0.94	ng		98
22) Phenanthrene	17.09	178	19463	1.16	ng		100
23) Anthracene	17.19	178	11922	0.90	ng		100
24) Fluoranthene	19.13	202	13640	1.01	ng		99
26) Pyrene	19.49	202	14483	1.00	ng		99
28) Benzo(a)anthracene	21.24	228	13512	1.00	ng		99
29) Chrysene	21.28	228	12782	1.02	ng		97
30) Bis(2-ethylhexyl)phthalate	21.16	149	7416	1.03	ng		99
31) Indeno(1,2,3-cd)pyrene	25.71	276	13282	1.00	ng		99
33) Benzo(b)fluoranthene	22.80	252	12309	1.00	ng		98
34) Benzo(k)fluoranthene	22.84	252	11904	1.03	ng		95
35) Benzo(a)pyrene	23.38	252	10832	1.01	ng		98
36) Dibenzo(a,h)anthracene	25.73	278	10749	1.01	ng		98
37) Benzo(g,h,i)perylene	26.40	276	11143	1.00	ng		99

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

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