Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
Data File : BM001532.D
Acq On : 03 Jun 2015 02:36
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001EC

Quant Time: Jun 03 03:36:33 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 03 02:19:00 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	17396	5.00	ng		0.00
6) Naphthalene-d8	10.49	136	61455	5.00	ng		0.00
12) Acenaphthene-d10	14.36	164	29797	5.00	ng		0.00
18) Phenanthrene-d10	17.09	188	91622	5.00	ng		0.00
25) Chrysene-d12	21.28	240	51143	5.00	ng		-0.01
32) Perylene-d12	23.53	264	48143	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.28	112	3872	1.02	ng		-0.02
5) Phenol-d6	6.86	99	4744	1.01	ng		0.00
7) Nitrobenzene-d5	8.87	82	3903	1.02	ng		0.00
13) 2,4,6-Tribromophenol	15.83	330	697	1.25	ng		0.00
14) 2-Fluorobiphenyl	12.98	172	8792	0.98	ng		0.00
27) Terphenyl-d14	19.73	244	6196	0.98	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.23	88	1626	0.99	ng		96
3) n-Nitrosodimethylamine	3.52	42	2556	1.03	ng		97
8) Nitrobenzene	8.91	77	4122	1.01	ng		98
9) Naphthalene	10.54	128	12747	0.98	ng		97
10) Hexachlorobutadiene	10.83	225	2259	1.00	ng		99
11) 2-Methylnaphthalene	12.17	142	8056	0.99	ng		91
15) Acenaphthylene	14.07	152	12447	0.98	ng		100
16) Acenaphthene	14.42	154	7854	0.99	ng		96
17) Fluorene	15.39	166	11146	0.84	ng		96
19) 4-Bromophenyl-phenylether	16.27	248	1211	0.95	ng	#	1
20) Hexachlorobenzene	16.41	284	2481	0.72	ng	#	37
21) Pentachlorophenol	16.75	266	1209	0.81	ng		92
22) Phenanthrene	17.12	178	18544	1.16	ng		99
23) Anthracene	17.23	178	14694	0.76	ng		99
24) Fluoranthene	19.16	202	14525	0.90	ng		98
26) Pyrene	19.53	202	15294	0.99	ng		100
28) Benzo(a)anthracene	21.27	228	13916	1.01	ng		98
29) Chrysene	21.31	228	12808	0.98	ng		100
30) Bis(2-ethylhexyl)phthalate	21.21	149	9284	0.95	ng	#	98
31) Indeno(1,2,3-cd)pyrene	25.79	276	14622	1.01	ng		94
33) Benzo(b)fluoranthene	22.85	252	13461	0.98	ng		96
34) Benzo(k)fluoranthene	22.90	252	12132	1.01	ng	#	96
35) Benzo(a)pyrene	23.43	252	11669	1.00	ng	#	96
36) Dibenzo(a,h)anthracene	25.82	278	11538	0.99	ng	#	95
37) Benzo(g,h,i)perylene	26.49	276	12097	0.99	ng		95

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

Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
Data File : BM001532.D
Acq On : 03 Jun 2015 02:36
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTDCCC001EC

Quant Time: Jun 03 03:36:33 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 03 02:19:00 2015
Response via : Initial Calibration