Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061915\
Data File : BM001752.D
Acq On : 19 Jun 2015 00:22
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTDCCC001EC

Manual Integrations APPROVED
apatel
6/22/2015 2:46:05 PM

Quant Time: Jun 19 05:17:37 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 05:05:14 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	15805	5.00	ng		0.00
6) Naphthalene-d8	10.44	136	57489	5.00	ng		-0.02
12) Acenaphthene-d10	14.31	164	27725	5.00	ng		-0.01
18) Phenanthrene-d10	17.06	188	67317	5.00	ng		0.00
25) Chrysene-d12	21.25	240	47893	5.00	ng		0.00
32) Perylene-d12	23.47	264	43059	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	3768	1.02	ng		0.00
5) Phenol-d6	6.83	99	4713	0.99	ng		0.00
7) Nitrobenzene-d5	8.83	82	4493	1.00	ng		0.00
13) 2,4,6-Tribromophenol	15.80	330	878	1.01	ng		0.00
14) 2-Fluorobiphenyl	12.94	172	9828	0.99	ng		0.00
27) Terphenyl-d14	19.70	244	6895	0.98	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.18	88	1827	1.01	ng		98
3) n-Nitrosodimethylamine	3.48	42	2797	1.01	ng	#	99
8) Nitrobenzene	8.87	77	4840	1.00	ng		99
9) Naphthalene	10.49	128	13896	0.99	ng		100
10) Hexachlorobutadiene	10.78	225	2338	0.99	ng		99
11) 2-Methylnaphthalene	12.12	142	8817	0.99	ng		99
15) Acenaphthylene	14.03	152	13106	0.99	ng		100
16) Acenaphthene	14.37	154	8188	0.99	ng		100
17) Fluorene	15.36	166	7636	1.08	ng		99
19) 4-Bromophenyl-phenylether	16.24	248	2811	1.02	ng		96
20) Hexachlorobenzene	16.38	284	1991	1.05	ng	#	100
21) Pentachlorophenol	16.72	266	944	1.12	ng		99
22) Phenanthrene	17.09	178	22432	0.97	ng		99
23) Anthracene	17.19	178	10038m	1.03	ng		
24) Fluoranthene	19.12	202	15450	1.06	ng		99
26) Pyrene	19.48	202	16269	0.98	ng		100
28) Benzo(a)anthracene	21.24	228	14199	0.98	ng		97
29) Chrysene	21.28	228	15105	1.00	ng		97
30) Bis(2-ethylhexyl)phthalate	21.16	149	8018	0.96	ng		100
31) Indeno(1,2,3-cd)pyrene	25.70	276	14586	1.00	ng		100
33) Benzo(b)fluoranthene	22.79	252	13252	0.94	ng		96
34) Benzo(k)fluoranthene	22.84	252	13382	1.06	ng		98
35) Benzo(a)pyrene	23.37	252	11848	1.00	ng		98
36) Dibenzo(a,h)anthracene	25.72	278	11782	1.00	ng		99
37) Benzo(g,h,i)perylene	26.39	276	12249	0.99	ng		99

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

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