Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
Data File : BM001534.D
Acq On : 03 Jun 2015 04:58
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Jun 03 05:42:16 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

Instrument :
BNA_M
ClientSampleId :
SSTDCCC001

Manual Integrations APPROVED
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Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.71	152	17305	5.00	ng		0.00
6) Naphthalene-d8	10.49	136	60932	5.00	ng		0.00
12) Acenaphthene-d10	14.36	164	29590	5.00	ng		0.00
18) Phenanthrene-d10	17.09	188	90106	5.00	ng		0.00
25) Chrysene-d12	21.28	240	49714	5.00	ng		-0.01
32) Perylene-d12	23.53	264	46672	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.28	112	3706	0.98	ng		-0.02
5) Phenol-d6	6.86	99	4525	0.97	ng		0.00
7) Nitrobenzene-d5	8.87	82	3780	1.00	ng		0.00
13) 2,4,6-Tribromophenol	15.83	330	685	1.24	ng		0.00
14) 2-Fluorobiphenyl	12.98	172	8829	0.99	ng		0.00
27) Terphenyl-d14	19.73	244	6136	1.00	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.23	88	1593	0.98	ng		99
3) n-Nitrosodimethylamine	3.52	42	2485	1.01	ng		100
8) Nitrobenzene	8.91	77	4019	1.00	ng		97
9) Naphthalene	10.54	128	12752	0.99	ng		96
10) Hexachlorobutadiene	10.83	225	2250	1.00	ng		100
11) 2-Methylnaphthalene	12.16	142	7973	0.99	ng		94
15) Acenaphthylene	14.07	152	12205	0.97	ng		100
16) Acenaphthene	14.42	154	7784	0.98	ng		97
17) Fluorene	15.39	166	10883	0.83	ng		98
19) 4-Bromophenyl-phenylether	16.27	248	1266	1.01	ng	#	1
20) Hexachlorobenzene	16.41	284	2523	0.75	ng	#	37
21) Pentachlorophenol	16.75	266	1215	0.82	ng		90
22) Phenanthrene	17.12	178	19051	1.21	ng		98
23) Anthracene	17.23	178	15014	0.79	ng		99
24) Fluoranthene	19.16	202	14204	0.90	ng		98
26) Pyrene	19.53	202	14929	1.00	ng		100
28) Benzo(a)anthracene	21.27	228	13663	1.02	ng		99
29) Chrysene	21.31	228	12107	0.96	ng		99
30) Bis(2-ethylhexyl)phthalate	21.21	149	8341	0.88	ng	#	97
31) Indeno(1,2,3-cd)pyrene	25.79	276	14085	1.00	ng		94
33) Benzo(b)fluoranthene	22.85	252	13183	0.99	ng		96
34) Benzo(k)fluoranthene	22.90	252	11613m	1.00	ng		
35) Benzo(a)pyrene	23.43	252	11118	0.98	ng	#	96
36) Dibenzo(a,h)anthracene	25.82	278	11139	0.99	ng		96
37) Benzo(g,h,i)perylene	26.48	276	11701	0.99	ng	#	94

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

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