Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061315\
Data File : BM001679.D
Acq On : 12 Jun 2015 20:55
Operator : TP/IZ
Sample : SSTDICC002
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDICC002

Quant Time: Jun 13 00:27:42 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	15474	5.00	ng		0.00
6) Naphthalene-d8	10.46	136	57054	5.00	ng		0.00
12) Acenaphthene-d10	14.32	164	28281	5.00	ng		0.00
18) Phenanthrene-d10	17.06	188	89014	5.00	ng		0.00
25) Chrysene-d12	21.25	240	48439	5.00	ng		0.00
32) Perylene-d12	23.48	264	43937	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	6670	1.81	ng		0.00
5) Phenol-d6	6.83	99	8536	1.84	ng		0.00
7) Nitrobenzene-d5	8.83	82	8060	1.83	ng		0.00
13) 2,4,6-Tribromophenol	15.80	330	1192	1.23	ng		0.00
14) 2-Fluorobiphenyl	12.94	172	17524	1.76	ng		0.00
27) Terphenyl-d14	19.70	244	13039	1.78	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.18	88	3169	1.63	ng		99
3) n-Nitrosodimethylamine	3.48	42	5020	1.85	ng		98
8) Nitrobenzene	8.87	77	8787	1.84	ng		99
9) Naphthalene	10.51	128	24303	1.75	ng		99
10) Hexachlorobutadiene	10.78	225	4048	1.74	ng		99
11) 2-Methylnaphthalene	12.13	142	15727	1.75	ng		100
15) Acenaphthylene	14.03	152	24657	1.85	ng		100
16) Acenaphthene	14.39	154	14797	1.75	ng		99
17) Fluorene	15.36	166	21375	2.33	ng		98
19) 4-Bromophenyl-phenylether	16.24	248	2961	0.38	ng		95
20) Hexachlorobenzene	16.38	284	5727	0.77	ng	#	100
21) Pentachlorophenol	16.72	266	2754	1.29	ng		98
22) Phenanthrene	17.09	178	36839	1.30	ng		98
23) Anthracene	17.19	178	28602	0.89	ng		100
24) Fluoranthene	19.13	202	28360	0.73	ng		100
26) Pyrene	19.49	202	29641	1.78	ng		99
28) Benzo(a)anthracene	21.24	228	27135	1.78	ng		99
29) Chrysene	21.28	228	25362	1.74	ng		99
30) Bis(2-ethylhexyl)phthalate	21.16	149	14717	1.56	ng		99
31) Indeno(1,2,3-cd)pyrene	25.72	276	26274	2.19	ng		100
33) Benzo(b)fluoranthene	22.80	252	23511	1.64	ng		98
34) Benzo(k)fluoranthene	22.85	252	24152	1.73	ng		98
35) Benzo(a)pyrene	23.38	252	21587	1.80	ng		98
36) Dibenzo(a,h)anthracene	25.73	278	21238	2.05	ng		99
37) Benzo(g,h,i)perylene	26.40	276	21786	1.95	ng		100

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

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