Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061315\
Data File : BM001685.D
Acq On : 13 Jun 2015 00:34
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM061315

Quant Time: Jun 13 02:15:23 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	14636	5.00	ng		0.00
6) Naphthalene-d8	10.46	136	54071	5.00	ng		0.00
12) Acenaphthene-d10	14.32	164	26415	5.00	ng		0.00
18) Phenanthrene-d10	17.06	188	81404	5.00	ng		0.00
25) Chrysene-d12	21.25	240	46120	5.00	ng		0.00
32) Perylene-d12	23.48	264	41923	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	3285	0.93	ng		0.00
5) Phenol-d6	6.83	99	4153	0.93	ng		0.00
7) Nitrobenzene-d5	8.83	82	3940	0.92	ng		0.00
13) 2,4,6-Tribromophenol	15.80	330	575	0.93	ng		0.00
14) 2-Fluorobiphenyl	12.94	172	8851	0.94	ng		0.00
27) Terphenyl-d14	19.70	244	6429	0.93	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.18	88	1613	1.02	ng		98
3) n-Nitrosodimethylamine	3.49	42	2517	0.95	ng		97
8) Nitrobenzene	8.87	77	4193	0.91	ng		100
9) Naphthalene	10.51	128	12202	0.94	ng		100
10) Hexachlorobutadiene	10.78	225	2073	0.95	ng		99
11) 2-Methylnaphthalene	12.13	142	7817	0.94	ng		100
15) Acenaphthylene	14.03	152	11709	0.93	ng		100
16) Acenaphthene	14.39	154	7161	0.92	ng		97
17) Fluorene	15.36	166	9695	0.85	ng		99
19) 4-Bromophenyl-phenylether	16.24	248	1514	0.98	ng		98
20) Hexachlorobenzene	16.38	284	2633	0.90	ng	#	100
21) Pentachlorophenol	16.72	266	1141	0.90	ng		99
22) Phenanthrene	17.09	178	18998	1.03	ng		100
23) Anthracene	17.19	178	13143	0.91	ng		99
24) Fluoranthene	19.13	202	13787	0.93	ng		100
26) Pyrene	19.49	202	14557	0.93	ng		99
28) Benzo(a)anthracene	21.24	228	13599	0.93	ng		98
29) Chrysene	21.28	228	12765	0.94	ng		97
30) Bis(2-ethylhexyl)phthalate	21.16	149	6811	0.92	ng		100
31) Indeno(1,2,3-cd)pyrene	25.71	276	13411	0.94	ng		100
33) Benzo(b)fluoranthene	22.80	252	12222	0.93	ng		98
34) Benzo(k)fluoranthene	22.84	252	11789	0.95	ng		96
35) Benzo(a)pyrene	23.38	252	10709	0.93	ng		98
36) Dibenzo(a,h)anthracene	25.73	278	10750	0.95	ng		98
37) Benzo(g,h,i)perylene	26.40	276	11262	0.94	ng		98

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

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