Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM052015\
Data File : BM001382.D
Acq On : 22 May 2015 04:52
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 22 05:36:25 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 25 10:49:04 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	114	0.00
2		1,4-Dioxane	1.000	0.614	38.6#	75	0.00
3		n-Nitrosodimethylamine	1.000	1.254	-25.4#	135	-0.02
4	S	2-Fluorophenol	1.000	1.860	-86.0#	216	0.00
5	S	Phenol-d6	1.000	1.784	-78.4#	206	0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	125	0.00
7	S	Nitrobenzene-d5	1.000	0.838	16.2	109	0.00
8		Nitrobenzene	1.000	0.807	19.3	110	0.00
9		Naphthalene	1.000	1.042	-4.2	131	0.00
10		Hexachlorobutadiene	1.000	0.951	4.9	120	0.00
11		2-Methylnaphthalene	1.000	1.076	-7.6	137	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	129	0.00
13	S	2,4,6-Tribromophenol	1.000	0.594	40.6#	74	0.00
14	S	2-Fluorobiphenyl	1.000	1.033	-3.3	134	0.00
15		Acenaphthylene	1.000	1.086	-8.6	145	0.00
16		Acenaphthene	1.000	1.049	-4.9	138	0.00
17		Fluorene	1.000	1.014	-1.4	134	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	131	0.00
19		4-Bromophenyl-phenylether	1.000	1.066	-6.6	145	0.00
20		Hexachlorobenzene	1.000	1.001	-0.1	132	0.00
21		Pentachlorophenol	1.000	0.739	26.1#	94	0.00
22		Phenanthrene	1.000	1.002	-0.2	130	0.00
23		Anthracene	1.000	0.994	0.6	122	0.00
24		Fluoranthene	1.000	1.055	-5.5	137	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	101	-0.02
26		Pyrene	1.000	1.284	-28.4#	131	0.00
27	S	Terphenyl-d14	1.000	1.294	-29.4#	133	0.00
28		Benzo(a)anthracene	1.000	1.104	-10.4	117	0.00
29		Chrysene	1.000	1.021	-2.1	106	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.584	-58.4#	205	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	0.797	20.3	87	0.00
32	I	Perylene-d12	5.000	5.000	0.0	83	0.00
33		Benzo(b)fluoranthene	1.000	1.034	-3.4	89	0.00
34		Benzo(k)fluoranthene	1.000	1.175	-17.5	99	-0.01
35	C	Benzo(a)pyrene	1.000	1.097	-9.7	96	0.00
36		Dibenzo(a,h)anthracene	1.000	1.000	0.0	88	0.00
37		Benzo(g,h,i)perylene	1.000	0.965	3.5	85	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 0