Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051815\
Data File : BM001331.D
Acq On : 18 May 2015 15:19
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 19 02:22:21 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 05:24:15 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	118	-0.02
2		1,4-Dioxane	1.000	0.749	25.1#	101	-0.01
3		n-Nitrosodimethylamine	1.000	1.050	-5.0	142	-0.02
4	S	2-Fluorophenol	1.000	0.928	7.2	116	-0.02
5	S	Phenol-d6	1.000	0.946	5.4	121	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	124	0.00
7	S	Nitrobenzene-d5	1.000	0.905	9.5	127	0.00
8		Nitrobenzene	1.000	0.891	10.9	128	0.00
9		Naphthalene	1.000	0.962	3.8	123	0.00
10		Hexachlorobutadiene	1.000	0.976	2.4	129	0.00
11		2-Methylnaphthalene	1.000	0.994	0.6	131	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	133	0.00
13	S	2,4,6-Tribromophenol	1.000	0.751	24.9	122	0.00
14	S	2-Fluorobiphenyl	1.000	0.944	5.6	132	0.00
15		Acenaphthylene	1.000	0.967	3.3	139	0.00
16		Acenaphthene	1.000	0.946	5.4	135	0.00
17		Fluorene	1.000	0.896	10.4	135	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	183	0.00
19		Hexachlorobenzene	1.000	0.698	30.2#	138	0.00
20		Pentachlorophenol	1.000	0.970	3.0	211	0.00
21		Phenanthrene	1.000	0.993	0.7	198	0.00
22		Anthracene	1.000	0.942	5.8	189	0.00
23		Fluoranthene	1.000	0.705	29.5#	143	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	134	0.00
25		Pyrene	1.000	0.994	0.6	146	0.00
26	S	Terphenyl-d14	1.000	0.991	0.9	141	-0.01
27		Benzo(a)anthracene	1.000	0.924	7.6	139	0.00
28		Chrysene	1.000	0.948	5.2	134	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	0.814	18.6	117	0.00
30	I	Perylene-d12	5.000	5.000	0.0	122	-0.01
31		Benzo(b)fluoranthene	1.000	1.003	-0.3	134	0.00
32		Benzo(k)fluoranthene	1.000	0.934	6.6	122	0.00
33	C	Benzo(a)pyrene	1.000	0.939	6.1	126	0.00
34		Dibenzo(a,h)anthracene	1.000	0.879	12.1	114	-0.01
35		Benzo(g,h,i)perylene	1.000	0.887	11.3	116	-0.01

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