Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071715\
Data File : BE090165.D
Acq On : 18 Jul 2015 2:59
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 18 04:52:29 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 00:40:53 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	82	0.00
2		1,4-Dioxane	1.000	1.116	-11.6	85	0.00
3		n-Nitrosodimethylamine	1.000	1.002	-0.2	84	0.00
4	S	2-Fluorophenol	1.000	0.911	8.9	75	0.00
5	S	Phenol-d6	1.000	0.940	6.0	76	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	80	0.00
7	S	Nitrobenzene-d5	1.000	0.933	6.7	77	0.00
8		Nitrobenzene	1.000	0.959	4.1	76	0.00
9		Naphthalene	1.000	0.983	1.7	81	0.00
10		Hexachlorobutadiene	1.000	0.970	3.0	79	0.00
11		2-Methylnaphthalene	1.000	0.951	4.9	78	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	76	0.00
13	S	2,4,6-Tribromophenol	1.000	0.752	24.8	59	0.00
14	S	2-Fluorobiphenyl	1.000	0.988	1.2	76	0.00
15		Acenaphthylene	1.000	0.973	2.7	76	0.00
16		Acenaphthene	1.000	0.970	3.0	76	0.00
17		Fluorene	1.000	0.950	5.0	74	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	76	0.00
19		4-Bromophenyl-phenylether	1.000	0.945	5.5	75	0.00
20		Hexachlorobenzene	1.000	0.945	5.5	74	0.00
21		Pentachlorophenol	1.000	0.751	24.9	55	0.00
22		Phenanthrene	1.000	0.968	3.2	76	0.00
23		Anthracene	1.000	0.941	5.9	74	0.02
24		Fluoranthene	1.000	0.955	4.5	75	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	77	0.00
26		Pyrene	1.000	0.956	4.4	75	0.00
27	S	Terphenyl-d14	1.000	0.952	4.8	75	0.00
28		Benzo(a)anthracene	1.000	0.921	7.9	74	0.00
29		Chrysene	1.000	0.981	1.9	78	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.713	28.7#	56	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.889	11.1	70	0.00
32	I	Perylene-d12	5.000	5.000	0.0	76	0.00
33		Benzo(b)fluoranthene	1.000	0.927	7.3	72	0.00
34		Benzo(k)fluoranthene	1.000	0.959	4.1	74	0.00
35	C	Benzo(a)pyrene	1.000	0.905	9.5	70	0.00
36		Dibenzo(a,h)anthracene	1.000	0.887	11.3	70	0.00
37		Benzo(g,h,i)perylene	1.000	0.919	8.1	71	0.00

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