Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071715\
Data File : BE090173.D
Acq On : 18 Jul 2015 9:00
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 20 04:58:58 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	93	0.00
2		1,4-Dioxane	1.000	1.099	-9.9	95	0.00
3		n-Nitrosodimethylamine	1.000	1.000	0.0	95	0.00
4	S	2-Fluorophenol	1.000	0.946	5.4	89	0.00
5	S	Phenol-d6	1.000	0.945	5.5	87	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
7	S	Nitrobenzene-d5	1.000	0.934	6.6	87	0.00
8		Nitrobenzene	1.000	0.961	3.9	87	0.00
9		Naphthalene	1.000	0.969	3.1	91	0.00
10		Hexachlorobutadiene	1.000	0.960	4.0	89	0.00
11		2-Methylnaphthalene	1.000	0.953	4.7	89	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	88	0.00
13	S	2,4,6-Tribromophenol	1.000	0.818	18.2	74	0.00
14	S	2-Fluorobiphenyl	1.000	0.983	1.7	88	0.00
15		Acenaphthylene	1.000	0.969	3.1	88	0.00
16		Acenaphthene	1.000	0.965	3.5	88	0.00
17		Fluorene	1.000	0.929	7.1	84	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	89	0.00
19		4-Bromophenyl-phenylether	1.000	0.947	5.3	87	0.00
20		Hexachlorobenzene	1.000	0.930	7.0	85	0.00
21		Pentachlorophenol	1.000	0.851	14.9	78	0.00
22		Phenanthrene	1.000	0.983	1.7	89	0.00
23		Anthracene	1.000	0.958	4.2	88	0.00
24		Fluoranthene	1.000	0.955	4.5	87	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	87	0.00
26		Pyrene	1.000	0.975	2.5	87	0.00
27	S	Terphenyl-d14	1.000	0.959	4.1	85	0.00
28		Benzo(a)anthracene	1.000	0.926	7.4	85	0.00
29		Chrysene	1.000	0.966	3.4	87	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.737	26.3#	66	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.862	13.8	77	0.00
32	I	Perylene-d12	5.000	5.000	0.0	84	0.00
33		Benzo(b)fluoranthene	1.000	0.933	6.7	80	0.00
34		Benzo(k)fluoranthene	1.000	0.961	3.9	83	0.00
35	C	Benzo(a)pyrene	1.000	0.912	8.8	78	0.00
36		Dibenzo(a,h)anthracene	1.000	0.884	11.6	77	0.00
37		Benzo(g,h,i)perylene	1.000	0.907	9.3	78	0.00

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