Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE063015\
Data File : BE090029.D
Acq On : 1 Jul 2015 2:55
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 01 06:20:16 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	96	0.00
2		1,4-Dioxane	1.000	0.859	14.1	86	-0.01
3		n-Nitrosodimethylamine	1.000	0.806	19.4	81	-0.01
4	S	2-Fluorophenol	1.000	0.965	3.5	93	0.00
5	S	Phenol-d6	1.000	0.965	3.5	94	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	96	-0.01
7	S	Nitrobenzene-d5	1.000	0.884	11.6	87	0.00
8		Nitrobenzene	1.000	0.864	13.6	86	0.00
9		Naphthalene	1.000	0.921	7.9	90	0.00
10		Hexachlorobutadiene	1.000	0.942	5.8	92	0.00
11		2-Methylnaphthalene	1.000	0.888	11.2	87	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	88	0.00
13	S	2,4,6-Tribromophenol	1.000	0.939	6.1	85	0.00
14	S	2-Fluorobiphenyl	1.000	0.923	7.7	84	0.00
15		Acenaphthylene	1.000	0.941	5.9	86	0.00
16		Acenaphthene	1.000	0.946	5.4	85	-0.01
17		Fluorene	1.000	0.902	9.8	82	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	81	-0.02
19		4-Bromophenyl-phenylether	1.000	0.879	12.1	73	0.00
20		Hexachlorobenzene	1.000	0.889	11.1	75	-0.02
21		Pentachlorophenol	1.000	0.883	11.7	73	0.00
22		Phenanthrene	1.000	0.884	11.6	74	0.00
23		Anthracene	1.000	0.932	6.8	77	0.00
24		Fluoranthene	1.000	0.907	9.3	75	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	81	-0.01
26		Pyrene	1.000	0.941	5.9	77	-0.01
27	S	Terphenyl-d14	1.000	0.901	9.9	73	-0.01
28		Benzo(a)anthracene	1.000	0.926	7.4	76	-0.01
29		Chrysene	1.000	0.898	10.2	73	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.098	-9.8	107	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	1.032	-3.2	85	-0.03
32	I	Perylene-d12	5.000	5.000	0.0	80	-0.02
33		Benzo(b)fluoranthene	1.000	1.024	-2.4	82	-0.01
34		Benzo(k)fluoranthene	1.000	0.995	0.5	80	-0.01
35	C	Benzo(a)pyrene	1.000	1.029	-2.9	82	-0.02
36		Dibenzo(a,h)anthracene	1.000	1.020	-2.0	83	-0.02
37		Benzo(g,h,i)perylene	1.000	1.027	-2.7	83	-0.03

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