Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE082515\
Data File : BE090627.D
Acq On : 25 Aug 2015 23:23
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 26 03:28:52 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	102	0.00
2		1,4-Dioxane	1.000	1.181	-18.1	103	0.00
3		n-Nitrosodimethylamine	1.000	1.016	-1.6	105	0.00
4	S	2-Fluorophenol	1.000	0.744	25.6#	75	0.00
5	S	Phenol-d6	1.000	0.678	32.2#	72	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
7	S	Nitrobenzene-d5	1.000	0.881	11.9	87	0.00
8		Nitrobenzene	1.000	0.885	11.5	89	0.00
9		Naphthalene	1.000	0.978	2.2	92	0.00
10		Hexachlorobutadiene	1.000	1.039	-3.9	98	0.00
11		2-Methylnaphthalene	1.000	0.927	7.3	89	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
13	S	2,4,6-Tribromophenol	1.000	0.858	14.2	95	0.00
14	S	2-Fluorobiphenyl	1.000	0.926	7.4	93	0.00
15		Acenaphthylene	1.000	0.938	6.2	96	0.00
16		Acenaphthene	1.000	0.982	1.8	100	0.00
17		Fluorene	1.000	1.028	-2.8	104	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	115	-0.02
19		4-Bromophenyl-phenylether	1.000	0.905	9.5	107	0.00
20		Hexachlorobenzene	1.000	0.973	2.7	113	-0.02
21		Pentachlorophenol	1.000	0.846	15.4	122	0.00
22		Phenanthrene	1.000	0.949	5.1	112	0.00
23		Anthracene	1.000	0.939	6.1	112	0.00
24		Fluoranthene	1.000	0.988	1.2	117	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	133	0.00
26		Pyrene	1.000	0.893	10.7	121	0.00
27	S	Terphenyl-d14	1.000	0.885	11.5	119	0.00
28		Benzo(a)anthracene	1.000	0.916	8.4	128	0.00
29		Chrysene	1.000	1.032	-3.2	138	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.679	32.1#	97	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.923	7.7	133	0.00
32	I	Perylene-d12	5.000	5.000	0.0	134	0.00
33		Benzo(b)fluoranthene	1.000	0.940	6.0	129	0.00
34		Benzo(k)fluoranthene	1.000	1.040	-4.0	138	0.00
35	C	Benzo(a)pyrene	1.000	0.946	5.4	131	0.00
36		Dibenzo(a,h)anthracene	1.000	0.911	8.9	129	0.00
37		Benzo(g,h,i)perylene	1.000	0.931	6.9	127	0.00

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