Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090335.D
Acq On : 4 Aug 2015 7:37
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 08:13:38 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	46#	0.00
2		1,4-Dioxane	0.796	0.799	-0.4	46#	0.00
3		n-Nitrosodimethylamine	0.845	0.894	-5.8	48#	0.00
4	S	2-Fluorophenol	0.956	1.011	-5.8	50	0.00
5	S	Phenol-d6	1.175	1.049	10.7	44#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	45#	0.00
7	S	Nitrobenzene-d5	0.295	0.288	2.4	46#	0.00
8		Nitrobenzene	0.305	0.273	10.5	43#	0.00
9		Naphthalene	1.074	1.077	-0.3	46#	0.00
10		Hexachlorobutadiene	0.144	0.143	0.7	44#	0.00
11		2-Methylnaphthalene	0.665	0.585	12.0	40#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	38#	0.00
13	S	2,4,6-Tribromophenol	0.067	0.054	19.4	42#	0.00
14	S	2-Fluorobiphenyl	1.449	1.545	-6.6	40#	0.00
15		Acenaphthylene	8.730	8.916	-2.1	39#	0.00
16		Acenaphthene	1.262	1.238	1.9	37#	0.00
17		Fluorene	1.533	1.444	5.8	38#	0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	34#	0.02
19		4-Bromophenyl-phenylether	0.188	0.214	-13.8	37#	0.00
20		Hexachlorobenzene	0.710	0.827	-16.5	38#	0.00
21		Pentachlorophenol	0.038	0.447	-1076.3#	613#	0.00
22		Phenanthrene	1.139	1.112	2.4	34#	0.00
23		Anthracene	1.018	0.984	3.3	35#	0.00
24		Fluoranthene	1.146	1.091	4.8	33#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	30#	0.00
26		Pyrene	1.352	1.590	-17.6	34#	0.00
27	S	Terphenyl-d14	0.936	1.062	-13.5	33#	0.00
28		Benzo(a)anthracene	0.951	0.885	6.9	31#	0.00
29		Chrysene	1.320	1.556	-17.9	34#	0.00
30		Bis(2-ethylhexyl)phthalate	0.335	0.284	15.2	36#	0.00
31		Indeno(1,2,3-cd)pyrene	0.548	0.702	-28.1#	54	0.00
32	I	Perylene-d12	1.000	1.000	0.0	30#	0.00
33		Benzo(b)fluoranthene	0.676	0.767	-13.5	39#	0.00
34		Benzo(k)fluoranthene	1.357	1.617	-19.2	33#	0.00
35	C	Benzo(a)pyrene	0.673	0.858	-27.5#	43#	0.00
36		Dibenzo(a,h)anthracene	0.431	0.615	-42.7#	61	0.00
37		Benzo(g,h,i)perylene	0.744	1.062	-42.7#	44#	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1