Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090333.D
Acq On : 4 Aug 2015 5:40
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 07:51:04 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	60	0.00
2		1,4-Dioxane	0.796	0.868	-9.0	65	-0.01
3		n-Nitrosodimethylamine	0.845	0.921	-9.0	65	0.00
4	S	2-Fluorophenol	0.956	1.151	-20.4	75	0.00
5	S	Phenol-d6	1.175	1.507	-28.3#	83	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	61	0.00
7	S	Nitrobenzene-d5	0.295	0.300	-1.7	65	0.00
8		Nitrobenzene	0.305	0.298	2.3	64	0.00
9		Naphthalene	1.074	1.194	-11.2	69	0.00
10		Hexachlorobutadiene	0.144	0.156	-8.3	65	0.00
11		2-Methylnaphthalene	0.665	0.741	-11.4	69	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	55	0.00
13	S	2,4,6-Tribromophenol	0.067	0.085	-26.9#	98	0.00
14	S	2-Fluorobiphenyl	1.449	1.812	-25.1#	68	0.00
15		Acenaphthylene	8.730	10.656	-22.1	67	0.00
16		Acenaphthene	1.262	1.479	-17.2	65	0.00
17		Fluorene	1.533	1.766	-15.2	67	0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	49#	0.02
19		4-Bromophenyl-phenylether	0.188	0.257	-36.7#	65	0.00
20		Hexachlorobenzene	0.710	0.945	-33.1#	64	0.00
21		Pentachlorophenol	0.038	0.044	-15.8	87	0.00
22		Phenanthrene	1.139	1.365	-19.8	61	0.00
23		Anthracene	1.018	1.235	-21.3	63	0.00
24		Fluoranthene	1.146	1.374	-19.9	61	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	45#	0.00
26		Pyrene	1.352	1.868	-38.2#	60	0.00
27	S	Terphenyl-d14	0.936	1.252	-33.8#	58	0.00
28		Benzo(a)anthracene	0.951	1.079	-13.5	58	0.00
29		Chrysene	1.320	1.942	-47.1#	64	0.00
30		Bis(2-ethylhexyl)phthalate	0.335	0.363	-8.4	69	0.00
31		Indeno(1,2,3-cd)pyrene	0.548	0.985	-79.7#	113	0.00
32	I	Perylene-d12	1.000	1.000	0.0	48#	0.00
33		Benzo(b)fluoranthene	0.676	1.116	-65.1#	92	0.00
34		Benzo(k)fluoranthene	1.357	2.151	-58.5#	71	0.00
35	C	Benzo(a)pyrene	0.673	1.111	-65.1#	90	0.00
36		Dibenzo(a,h)anthracene	0.431	0.706	-63.8#	112	0.00
37		Benzo(g,h,i)perylene	0.744	1.403	-88.6#	93	-0.01

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