Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090342.D
Acq On : 4 Aug 2015 11:50
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 13:22:23 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	47	0.00
2		1,4-Dioxane	1.000	2.006	-100.6#	73	0.00
3		n-Nitrosodimethylamine	1.000	1.043	-4.3	49	0.00
4	S	2-Fluorophenol	1.000	1.411	-41.1#	69	0.00
5	S	Phenol-d6	1.000	1.307	-30.7#	67	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	53	0.00
7	S	Nitrobenzene-d5	1.000	1.461	-46.1#	80	0.00
8		Nitrobenzene	1.000	0.806	19.4	45	0.00
9		Naphthalene	1.000	0.824	17.6	44	0.00
10		Hexachlorobutadiene	1.000	0.841	15.9	43	0.00
11		2-Methylnaphthalene	1.000	0.756	24.4	40	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	44	0.00
13	S	2,4,6-Tribromophenol	1.000	2.003	-100.3#	121	0.00
14	S	2-Fluorobiphenyl	1.000	1.769	-76.9#	77	0.00
15		Acenaphthylene	1.000	0.820	18.0	36	0.00
16		Acenaphthene	1.000	0.769	23.1	34	0.00
17		Fluorene	1.000	0.732	26.8#	34	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	39	0.00
19		4-Bromophenyl-phenylether	1.000	0.873	12.7	33	0.00
20		Hexachlorobenzene	1.000	0.872	12.8	33	0.00
21		Pentachlorophenol	1.000	1.103	-10.3	53	0.00
22		Phenanthrene	1.000	0.777	22.3	31	0.00
23		Anthracene	1.000	0.743	25.7#	30	0.00
24		Fluoranthene	1.000	0.758	24.2	31	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	36	0.00
26		Pyrene	1.000	0.866	13.4	30	0.00
27	S	Terphenyl-d14	1.000	1.917	-91.7#	66	0.00
28		Benzo(a)anthracene	1.000	0.695	30.5#	28	0.00
29		Chrysene	1.000	0.872	12.8	27	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.995	0.5	43	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.176	-17.6	56	0.00
32	I	Perylene-d12	5.000	5.000	0.0	40	0.00
33		Benzo(b)fluoranthene	1.000	0.948	5.2	43	0.00
34		Benzo(k)fluoranthene	1.000	0.717	28.3#	27	0.00
35	C	Benzo(a)pyrene	1.000	0.946	5.4	41	0.00
36		Dibenzo(a,h)anthracene	1.000	0.928	7.2	47	0.00
37		Benzo(g,h,i)perylene	1.000	0.996	0.4	41	-0.01

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