Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082615\
Data File : BE090635.D
Acq On : 26 Aug 2015 11:52
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 26 23:41:22 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	109	0.00
2		1,4-Dioxane	0.658	0.700	-6.4	109	0.00
3		n-Nitrosodimethylamine	0.863	0.880	-2.0	112	0.00
4	S	2-Fluorophenol	1.446	1.060	26.7#	79	0.00
5	S	Phenol-d6	2.125	1.469	30.9#	78	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
7	S	Nitrobenzene-d5	0.398	0.373	6.3	99	0.00
8		Nitrobenzene	0.413	0.373	9.7	97	0.00
9		Naphthalene	1.136	1.106	2.6	98	0.00
10		Hexachlorobutadiene	0.167	0.173	-3.6	104	0.00
11		2-Methylnaphthalene	0.740	0.687	7.2	95	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
13	S	2,4,6-Tribromophenol	0.192	0.155	19.3	97	0.00
14	S	2-Fluorobiphenyl	1.493	1.348	9.7	99	0.00
15		Acenaphthylene	9.132	8.472	7.2	104	0.00
16		Acenaphthene	1.374	1.342	2.3	108	0.00
17		Fluorene	1.684	1.767	-4.9	116	0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	127	0.00
19		4-Bromophenyl-phenylether	0.204	0.179	12.3	114	0.00
20		Hexachlorobenzene	0.906	0.874	3.5	123	-0.02
21		Pentachlorophenol	0.079	0.056	29.1#	119	0.00
22		Phenanthrene	1.312	1.231	6.2	121	0.00
23		Anthracene	1.208	1.115	7.7	121	0.00
24		Fluoranthene	1.396	1.357	2.8	126	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	143	0.00
26		Pyrene	1.430	1.287	10.0	132	0.00
27	S	Terphenyl-d14	0.914	0.799	12.6	127	0.00
28		Benzo(a)anthracene	1.316	1.203	8.6	138	0.00
29		Chrysene	1.344	1.385	-3.1	149	0.00
30		Bis(2-ethylhexyl)phthalate	0.725	0.423	41.7#	102	0.00
31		Indeno(1,2,3-cd)pyrene	1.266	1.156	8.7	142	0.00
32	I	Perylene-d12	1.000	1.000	0.0	143	0.00
33		Benzo(b)fluoranthene	1.184	1.173	0.9	146	0.00
34		Benzo(k)fluoranthene	1.342	1.354	-0.9	143	0.00
35	C	Benzo(a)pyrene	1.108	1.048	5.4	140	0.00
36		Dibenzo(a,h)anthracene	1.069	0.977	8.6	138	0.00
37		Benzo(g,h,i)perylene	1.169	1.084	7.3	135	0.00

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