Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE040417\
Data File : BE092886.D
Acq On : 4 Apr 2017 15:58
Operator : SJ/MA
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE040417

Quant Time: Apr 04 16:37:50 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 04 16:34:01 2017
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	107	-0.02
2		1,4-Dioxane	0.671	0.684	-1.9	109	0.00
3		n-Nitrosodimethylamine	0.673	0.663	1.5	112	0.00
4	S	2-Fluorophenol	1.007	1.007	0.0	117	0.00
5	S	Phenol-d6	1.445	1.396	3.4	117	0.00
6		bis(2-Chloroethyl)ether	1.622	1.584	2.3	107	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
8	S	Nitrobenzene-d5	0.287	0.298	-3.8	126	-0.02
9		Naphthalene	1.103	1.106	-0.3	109	0.00
10		Hexachlorobutadiene	0.149	0.148	0.7	108	0.00
11		2-Methylnaphthalene	0.684	0.672	1.8	110	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	113	-0.02
13	S	2,4,6-Tribromophenol	0.064	0.059	7.8	123	-0.01
14	S	2-Fluorobiphenyl	1.689	1.663	1.5	109	-0.01
15		Acenaphthylene	7.071	6.531	7.6	117	-0.02
16		Acenaphthene	1.452	1.398	3.7	112	0.00
17		Fluorene	1.606	1.517	5.5	112	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	113	-0.02
19		Hexachlorobenzene	0.190	0.182	4.2	109	0.00
20		Pentachlorophenol	0.036	0.034	5.6	116	-0.02
21		Phenanthrene	1.335	1.297	2.8	113	0.00
22		Anthracene	1.090	0.973	10.7	114	0.00
23		Fluoranthene	4.886	4.496	8.0	111	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	110	-0.02
25		Pyrene	4.464	4.245	4.9	113	0.00
26	S	Terphenyl-d14	0.768	0.739	3.8	111	-0.02
27		Benzo(a)anthracene	1.079	0.951	11.9	114	0.00
28		Chrysene	1.252	1.283	-2.5	108	0.00
29		Bis(2-ethylhexyl)phthalate	0.260	0.223	14.2	121	-0.02
30		Indeno(1,2,3-cd)pyrene	4.172	3.710	11.1	101	-0.02
31	I	Perylene-d12	1.000	1.000	0.0	105	-0.01
32		Benzo(b)fluoranthene	1.236	1.140	7.8	104	-0.01
33		Benzo(k)fluoranthene	1.119	1.014	9.4	107	-0.01
34	C	Benzo(a)pyrene	0.961	0.831	13.5	104	-0.01
35		Dibenzo(a,h)anthracene	1.102	0.980	11.1	102	-0.03
36		Benzo(g,h,i)perylene	4.504	4.128	8.3	102	-0.02

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