Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE110816\
Data File : BE092509.D
Acq On : 8 Nov 2016 15:16
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE110816

Quant Time: Nov 08 16:31:51 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE110816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 08 16:27:22 2016
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	105	0.00
2		1,4-Dioxane	0.775	0.826	-6.6	127	0.00
3		n-Nitrosodimethylamine	0.905	0.751	17.0	100	0.00
4	S	2-Fluorophenol	1.103	1.052	4.6	109	0.00
5	S	Phenol-d6	1.471	1.315	10.6	107	0.00
6		bis(2-Chloroethyl)ether	1.610	1.465	9.0	103	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
8	S	Nitrobenzene-d5	0.328	0.304	7.3	111	0.00
9		Naphthalene	1.082	1.070	1.1	102	0.00
10		Hexachlorobutadiene	0.164	0.169	-3.0	104	-0.02
11		2-Methylnaphthalene	0.703	0.687	2.3	106	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
13	S	2,4,6-Tribromophenol	0.142	0.111	21.8	106	0.00
14	S	2-Fluorobiphenyl	1.237	1.263	-2.1	111	0.00
15		Acenaphthylene	7.172	6.975	2.7	109	-0.02
16		Acenaphthene	1.159	1.143	1.4	107	0.00
17		Fluorene	1.455	1.422	2.3	108	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
19		Hexachlorobenzene	0.189	0.193	-2.1	112	0.00
20		Pentachlorophenol	0.055	0.044	20.0	97	0.00
21		Phenanthrene	1.081	1.156	-6.9	113	0.00
22		Anthracene	1.048	1.048	0.0	107	0.00
23		Fluoranthene	5.427	5.455	-0.5	106	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	108	0.00
25		Pyrene	4.097	3.757	8.3	107	0.00
26	S	Terphenyl-d14	0.607	0.566	6.8	109	-0.02
27		Benzo(a)anthracene	5.068	4.809	5.1	111	0.00
28		Chrysene	5.430	5.192	4.4	101	0.00
29		Bis(2-ethylhexyl)phthalate	0.535	0.445	16.8	102	0.00
30		Indeno(1,2,3-cd)pyrene	5.531	5.280	4.5	102	0.00
31	I	Perylene-d12	1.000	1.000	0.0	103	0.00
32		Benzo(b)fluoranthene	1.885	1.959	-3.9	103	0.00
33		Benzo(k)fluoranthene	1.972	1.824	7.5	97	0.00
34	C	Benzo(a)pyrene	1.292	1.233	4.6	102	0.00
35		Dibenzo(a,h)anthracene	1.216	1.174	3.5	102	-0.02
36		Benzo(g,h,i)perylene	5.009	4.857	3.0	103	0.00

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