Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082916\
Data File : BE092293.D
Acq On : 29 Aug 2016 17:37
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE082916

Quant Time: Aug 29 18:43:42 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE082916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 29 18:19:23 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	101	0.00
2		1,4-Dioxane	0.711	0.788	-10.8	108	0.00
3		n-Nitrosodimethylamine	0.740	0.781	-5.5	122	-0.02
4	S	2-Fluorophenol	1.122	1.111	1.0	119	-0.02
5	S	Phenol-d6	1.478	1.286	13.0	111	-0.02
6		bis(2-Chloroethyl)ether	1.151	1.230	-6.9	128	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	105	-0.02
8	S	Nitrobenzene-d5	0.327	0.273	16.5	117	0.00
9		Naphthalene	1.115	1.095	1.8	103	0.00
10		Hexachlorobutadiene	0.169	0.170	-0.6	105	0.00
11		2-Methylnaphthalene	0.689	0.669	2.9	107	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
13	S	2,4,6-Tribromophenol	0.135	0.122	9.6	106	0.00
14	S	2-Fluorobiphenyl	1.570	1.542	1.8	108	0.00
15		Acenaphthylene	8.560	8.110	5.3	107	0.00
16		Acenaphthene	1.397	1.362	2.5	107	0.00
17		Fluorene	1.689	1.642	2.8	108	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
19		Hexachlorobenzene	0.206	0.194	5.8	106	0.00
20		Pentachlorophenol	0.049	0.039	20.4	118	0.00
21		Phenanthrene	1.177	1.161	1.4	110	0.00
22		Anthracene	1.080	1.021	5.5	108	0.00
23		Fluoranthene	5.268	4.793	9.0	105	-0.02
24	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
25		Pyrene	4.627	4.410	4.7	108	0.00
26	S	Terphenyl-d14	0.676	0.647	4.3	104	0.00
27		Benzo(a)anthracene	5.171	4.451	13.9	107	0.00
28		Chrysene	4.699	4.916	-4.6	104	0.00
29		Bis(2-ethylhexyl)phthalate	0.328	0.289	11.9	103	0.00
30		Indeno(1,2,3-cd)pyrene	4.773	4.565	4.4	111	0.00
31	I	Perylene-d12	1.000	1.000	0.0	104	0.00
32		Benzo(b)fluoranthene	1.185	1.150	3.0	102	0.00
33		Benzo(k)fluoranthene	1.343	1.293	3.7	107	0.00
34	C	Benzo(a)pyrene	1.186	1.108	6.6	105	0.01
35		Dibenzo(a,h)anthracene	1.040	0.985	5.3	113	0.00
36		Benzo(g,h,i)perylene	4.652	4.489	3.5	108	0.00

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