Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042017\
Data File : BE092904.D
Acq On : 20 Apr 2017 17:35
Operator : SJ/MA
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE042017

Quant Time: Apr 20 18:07:31 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 20 17:35:10 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	94	0.00
2		1,4-Dioxane	0.768	0.740	3.6	97	0.00
3		n-Nitrosodimethylamine	0.582	0.576	1.0	98	0.00
4	S	2-Fluorophenol	0.870	0.837	3.8	98	0.00
5	S	Phenol-d6	1.213	1.171	3.5	100	0.00
6		bis(2-Chloroethyl)ether	1.474	1.410	4.3	95	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
8	S	Nitrobenzene-d5	0.267	0.262	1.9	96	0.00
9		Naphthalene	1.066	1.039	2.5	94	0.00
10		Hexachlorobutadiene	0.148	0.152	-2.7	98	0.00
11	SURR	2-Methylnaphthalene-d10	0.590	0.586	0.7	97	0.00
12		2-Methylnaphthalene	0.658	0.641	2.6	96	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
14	S	2,4,6-Tribromophenol	0.066	0.056	15.2	93	0.00
15	S	2-Fluorobiphenyl	1.838	1.866	-1.5	95	0.00
16		Acenaphthylene	6.430	6.064	5.7	95	0.00
17		Acenaphthene	1.263	1.223	3.2	93	0.00
18		Fluorene	1.541	1.504	2.4	95	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
20		Hexachlorobenzene	0.177	0.161	9.0	89	0.00
21		Pentachlorophenol	0.032	0.028	12.5	99	0.00
22		Phenanthrene	1.159	1.120	3.4	94	0.00
23		Anthracene	0.886	0.819	7.6	93	0.00
24	SURR	Fluoranthene-d10	0.958	0.902	5.8	96	0.00
25		Fluoranthene	4.994	4.867	2.5	100	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	92	0.00
27		Pyrene	5.224	5.221	0.1	96	0.00
28	S	Terphenyl-d14	0.694	0.688	0.9	92	0.00
29		Benzo(a)anthracene	0.931	0.890	4.4	97	-0.02
30		Chrysene	1.193	1.212	-1.6	99	0.00
31		Bis(2-ethylhexyl)phthalate	0.241	0.244	-1.2	103	0.00
32		Indeno(1,2,3-cd)pyrene	3.910	4.001	-2.3	103	0.00
33	I	Perylene-d12	1.000	1.000	0.0	98	0.00
34		Benzo(b)fluoranthene	1.401	1.379	1.6	100	0.00
35		Benzo(k)fluoranthene	1.427	1.344	5.8	99	0.00
36	C	Benzo(a)pyrene	1.133	1.062	6.3	99	0.00
37		Dibenzo(a,h)anthracene	1.264	1.252	0.9	105	0.00
38		Benzo(g,h,i)perylene	5.484	5.529	-0.8	103	0.02

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