Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE092216\
Data File : BE092400.D
Acq On : 22 Sep 2016 17:49
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE092216

Quant Time: Sep 22 18:39:58 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Sep 22 18:25:56 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	103	-0.02
2		1,4-Dioxane	0.635	0.607	4.4	97	-0.01
3		n-Nitrosodimethylamine	0.588	0.512	12.9	99	-0.01
4	S	2-Fluorophenol	0.999	0.833	16.6	92	0.00
5	S	Phenol-d6	1.388	1.048	24.5	89	0.00
6		bis(2-Chloroethyl)ether	1.222	0.839	31.3#	91	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
8	S	Nitrobenzene-d5	0.299	0.237	20.7	94	0.00
9		Naphthalene	1.082	1.112	-2.8	103	0.00
10		Hexachlorobutadiene	0.170	0.171	-0.6	101	0.00
11		2-Methylnaphthalene	0.716	0.716	0.0	101	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	0.144	0.114	20.8	95	0.01
14	S	2-Fluorobiphenyl	1.184	1.107	6.5	97	0.01
15		Acenaphthylene	7.240	6.880	5.0	100	0.02
16		Acenaphthene	1.134	1.156	-1.9	101	0.00
17		Fluorene	1.438	1.376	4.3	100	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.02
19		Hexachlorobenzene	0.214	0.197	7.9	99	0.00
20		Pentachlorophenol	0.050	0.041	18.0	97	0.00
21		Phenanthrene	1.149	1.083	5.7	100	0.02
22		Anthracene	1.105	0.998	9.7	100	0.02
23		Fluoranthene	5.460	5.053	7.5	101	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
25		Pyrene	4.333	4.107	5.2	99	0.02
26	S	Terphenyl-d14	0.620	0.583	6.0	96	0.02
27		Benzo(a)anthracene	4.873	4.612	5.4	97	0.00
28		Chrysene	4.462	4.654	-4.3	98	0.00
29		Bis(2-ethylhexyl)phthalate	0.420	0.337	19.8	80	0.00
30		Indeno(1,2,3-cd)pyrene	4.734	4.314	8.9	93	0.05
31	I	Perylene-d12	1.000	1.000	0.0	98	0.03
32		Benzo(b)fluoranthene	1.211	1.208	0.2	98	0.03
33		Benzo(k)fluoranthene	1.267	1.180	6.9	98	0.03
34	C	Benzo(a)pyrene	1.187	1.124	5.3	98	0.02
35		Dibenzo(a,h)anthracene	1.088	1.021	6.2	96	0.05
36		Benzo(g,h,i)perylene	4.673	4.526	3.1	98	0.05

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