Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091815\
Data File : BE090890.D
Acq On : 18 Sep 2015 17:49
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE091815

Quant Time: Sep 18 18:33:10 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 18 18:17:34 2015
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	102	0.00
2		1,4-Dioxane	1.000	0.991	0.9	93	0.00
3		n-Nitrosodimethylamine	1.000	0.964	3.6	92	0.00
4	S	2-Fluorophenol	1.000	0.918	8.2	90	0.00
5	S	Phenol-d6	1.000	0.886	11.4	93	0.00
6		bis(2-Chloroethyl)ether	1.000	1.015	-1.5	105	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	101	0.00
8	S	Nitrobenzene-d5	1.000	0.920	8.0	94	0.00
9		Nitrobenzene	1.000	0.907	9.3	95	0.00
10		Naphthalene	1.000	0.976	2.4	94	0.00
11		Hexachlorobutadiene	1.000	0.982	1.8	95	0.00
12		2-Methylnaphthalene	1.000	0.918	8.2	95	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
14	S	2,4,6-Tribromophenol	1.000	0.853	14.7	90	-0.02
15	S	2-Fluorobiphenyl	1.000	0.939	6.1	96	0.00
16		Acenaphthylene	1.000	0.916	8.4	95	0.00
17		Acenaphthene	1.000	0.930	7.0	97	0.00
18		Fluorene	1.000	0.952	4.8	97	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
20		4-Bromophenyl-phenylether	1.000	0.930	7.0	96	-0.02
21		Hexachlorobenzene	1.000	0.987	1.3	96	0.00
22		Pentachlorophenol	1.000	0.800	20.0	77	0.00
23		Phenanthrene	1.000	1.009	-0.9	99	0.00
24		Anthracene	1.000	0.901	9.9	95	0.00
25		Fluoranthene	1.000	0.898	10.2	93	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	95	0.00
27		Pyrene	1.000	0.903	9.7	92	0.00
28	S	Terphenyl-d14	1.000	0.920	8.0	91	0.00
29		Benzo(a)anthracene	1.000	0.875	12.5	88	0.00
30		Chrysene	1.000	0.965	3.5	88	0.00
31		Bis(2-ethylhexyl)phthalate	1.000	0.768	23.2	75	0.00
32		Indeno(1,2,3-cd)pyrene	1.000	0.871	12.9	83	0.00
33	I	Perylene-d12	5.000	5.000	0.0	92	0.00
34		Benzo(b)fluoranthene	1.000	0.868	13.2	83	0.00
35		Benzo(k)fluoranthene	1.000	0.923	7.7	86	0.00
36	C	Benzo(a)pyrene	1.000	0.866	13.4	83	0.00
37		Dibenzo(a,h)anthracene	1.000	0.872	12.8	81	0.00
38		Benzo(g,h,i)perylene	1.000	0.907	9.3	85	0.00

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