Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
Data File : BE090523.D
Acq On : 14 Aug 2015 16:43
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE081415

Quant Time: Aug 17 04:17:49 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	136	0.00
2		1,4-Dioxane	1.000	1.066	-6.6	125	0.00
3		n-Nitrosodimethylamine	1.000	0.962	3.8	133	0.00
4	S	2-Fluorophenol	1.000	0.947	5.3	128	0.00
5	S	Phenol-d6	1.000	0.962	3.8	137	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	135	0.00
7	S	Nitrobenzene-d5	1.000	0.971	2.9	138	0.00
8		Nitrobenzene	1.000	0.967	3.3	140	0.00
9		Naphthalene	1.000	0.965	3.5	131	0.00
10		Hexachlorobutadiene	1.000	0.988	1.2	134	0.00
11		2-Methylnaphthalene	1.000	0.967	3.3	134	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	137	0.00
13	S	2,4,6-Tribromophenol	1.000	0.917	8.3	139	-0.02
14	S	2-Fluorobiphenyl	1.000	0.975	2.5	134	0.00
15		Acenaphthylene	1.000	0.971	2.9	136	0.00
16		Acenaphthene	1.000	0.975	2.5	135	0.00
17		Fluorene	1.000	0.991	0.9	137	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	137	0.00
19		4-Bromophenyl-phenylether	1.000	0.973	2.7	137	0.00
20		Hexachlorobenzene	1.000	0.968	3.2	133	-0.02
21		Pentachlorophenol	1.000	0.881	11.9	154	0.00
22		Phenanthrene	1.000	0.937	6.3	131	0.00
23		Anthracene	1.000	0.965	3.5	137	0.00
24		Fluoranthene	1.000	0.953	4.7	134	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	138	0.00
26		Pyrene	1.000	0.968	3.2	136	0.00
27	S	Terphenyl-d14	1.000	0.976	2.4	136	0.00
28		Benzo(a)anthracene	1.000	0.945	5.5	137	0.00
29		Chrysene	1.000	0.969	3.1	134	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.903	9.7	142	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.983	1.7	146	0.00
32	I	Perylene-d12	5.000	5.000	0.0	139	0.00
33		Benzo(b)fluoranthene	1.000	0.956	4.4	136	0.00
34		Benzo(k)fluoranthene	1.000	1.019	-1.9	141	0.00
35	C	Benzo(a)pyrene	1.000	0.977	2.3	140	0.00
36		Dibenzo(a,h)anthracene	1.000	0.998	0.2	146	0.00
37		Benzo(g,h,i)perylene	1.000	0.987	1.3	140	0.00

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