Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071615\
Data File : BE090133.D
Acq On : 16 Jul 2015 23:54
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE071615

Quant Time: Jul 17 13:00:54 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 17 12:59:41 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	87	0.00
2		1,4-Dioxane	1.000	1.060	-6.0	90	0.00
3		n-Nitrosodimethylamine	1.000	0.950	5.0	87	0.00
4	S	2-Fluorophenol	1.000	0.996	0.4	93	0.00
5	S	Phenol-d6	1.000	1.021	-2.1	93	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	88	0.00
7	S	Nitrobenzene-d5	1.000	0.983	1.7	90	0.00
8		Nitrobenzene	1.000	0.993	0.7	91	0.00
9		Naphthalene	1.000	0.971	2.9	89	0.00
10		Hexachlorobutadiene	1.000	0.958	4.2	88	0.00
11		2-Methylnaphthalene	1.000	0.994	0.6	91	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
13	S	2,4,6-Tribromophenol	1.000	0.919	8.1	106	0.00
14	S	2-Fluorobiphenyl	1.000	0.971	2.9	89	0.00
15		Acenaphthylene	1.000	0.980	2.0	91	0.00
16		Acenaphthene	1.000	0.967	3.3	90	0.00
17		Fluorene	1.000	0.963	3.7	89	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	92	0.00
19		4-Bromophenyl-phenylether	1.000	0.975	2.5	93	0.00
20		Hexachlorobenzene	1.000	0.963	3.7	89	0.00
21		Pentachlorophenol	1.000	1.038	-3.8	102	0.00
22		Phenanthrene	1.000	0.955	4.5	91	0.00
23		Anthracene	1.000	0.965	3.5	91	0.00
24		Fluoranthene	1.000	0.996	0.4	93	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	97	0.00
26		Pyrene	1.000	0.928	7.2	92	0.00
27	S	Terphenyl-d14	1.000	0.989	1.1	95	0.00
28		Benzo(a)anthracene	1.000	1.070	-7.0	110	0.00
29		Chrysene	1.000	1.036	-3.6	97	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.979	2.1	101	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.178	-17.8	123	0.00
32	I	Perylene-d12	5.000	5.000	0.0	98	0.00
33		Benzo(b)fluoranthene	1.000	1.324	-32.4#	132	0.00
34		Benzo(k)fluoranthene	1.000	0.947	5.3	104	0.00
35	C	Benzo(a)pyrene	1.000	0.942	5.8	124	0.00
36		Dibenzo(a,h)anthracene	1.000	1.247	-24.7	133	0.00
37		Benzo(g,h,i)perylene	1.000	0.941	5.9	117	0.00

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