Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051315\
Data File : BM001267.D
Acq On : 13 May 2015 00:54
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM051315

Quant Time: May 13 05:49:29 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 05:24:15 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	93	0.00
2		1,4-Dioxane	1.000	1.084	-8.4	98	0.00
3		n-Nitrosodimethylamine	1.000	1.003	-0.3	107	0.00
4	S	2-Fluorophenol	1.000	0.947	5.3	93	0.00
5	S	Phenol-d6	1.000	0.917	8.3	92	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
7	S	Nitrobenzene-d5	1.000	0.868	13.2	90	0.00
8		Nitrobenzene	1.000	0.872	12.8	93	0.00
9		Naphthalene	1.000	0.970	3.0	92	0.00
10		Hexachlorobutadiene	1.000	0.954	4.6	93	0.00
11		2-Methylnaphthalene	1.000	0.942	5.8	92	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	91	0.00
13	S	2,4,6-Tribromophenol	1.000	0.726	27.4#	80	0.00
14	S	2-Fluorobiphenyl	1.000	0.968	3.2	93	0.00
15		Acenaphthylene	1.000	0.926	7.4	91	0.00
16		Acenaphthene	1.000	0.946	5.4	92	0.00
17		Fluorene	1.000	0.859	14.1	88	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	105	0.00
19		Hexachlorobenzene	1.000	0.845	15.5	96	0.00
20		Pentachlorophenol	1.000	0.842	15.8	93	0.00
21		Phenanthrene	1.000	0.845	15.5	97	0.00
22		Anthracene	1.000	0.829	17.1	95	0.00
23		Fluoranthene	1.000	0.788	21.2	92	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	92	-0.01
25		Pyrene	1.000	0.925	7.5	93	0.00
26	S	Terphenyl-d14	1.000	0.926	7.4	91	-0.01
27		Benzo(a)anthracene	1.000	0.902	9.8	94	0.00
28		Chrysene	1.000	0.961	3.9	94	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	0.977	2.3	96	0.00
30	I	Perylene-d12	5.000	5.000	0.0	93	-0.01
31		Benzo(b)fluoranthene	1.000	0.961	3.9	98	0.00
32		Benzo(k)fluoranthene	1.000	0.901	9.9	90	0.00
33	C	Benzo(a)pyrene	1.000	0.909	9.1	93	0.00
34		Dibenzo(a,h)anthracene	1.000	0.962	3.8	96	-0.01
35		Benzo(g,h,i)perylene	1.000	0.973	2.7	97	-0.01

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