Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120215\
Data File : BM003359.D
Acq On : 02 Dec 2015 16:57
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM120215

Quant Time: Dec 02 17:37:21 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 17:30:44 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	68	0.00
2		1,4-Dioxane	1.000	0.929	7.1	71	0.00
3		n-Nitrosodimethylamine	1.000	0.955	4.5	67	0.00
4	S	2-Fluorophenol	1.000	0.908	9.2	69	0.00
5	S	Phenol-d6	1.000	0.919	8.1	66	0.00
6		bis(2-Chloroethyl)ether	1.000	0.953	4.7	67	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	66	0.00
8	S	Nitrobenzene-d5	1.000	0.922	7.8	66	0.00
9		Nitrobenzene	1.000	0.921	7.9	64	0.00
10		Naphthalene	1.000	0.946	5.4	68	0.00
11		Hexachlorobutadiene	1.000	0.982	1.8	70	0.00
12		2-Methylnaphthalene	1.000	0.943	5.7	66	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	64	0.00
14	S	2,4,6-Tribromophenol	1.000	0.937	6.3	59	0.00
15	S	2-Fluorobiphenyl	1.000	0.922	7.8	66	0.00
16		Acenaphthylene	1.000	0.937	6.3	62	0.00
17		Acenaphthene	1.000	0.896	10.4	65	0.00
18		Fluorene	1.000	0.907	9.3	62	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	64	0.00
20		4-Bromophenyl-phenylether	1.000	0.852	14.8	63	0.00
21		Hexachlorobenzene	1.000	0.915	8.5	64	0.00
22		Pentachlorophenol	1.000	0.848	15.2	52	0.00
23		Phenanthrene	1.000	0.898	10.2	64	0.00
24		Anthracene	1.000	0.897	10.3	58	0.00
25		Fluoranthene	1.000	0.903	9.7	57	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	51	0.00
27		Benzidine	1.000	0.930	7.0	52	0.02
28		Pyrene	1.000	0.979	2.1	55	0.00
29	S	Terphenyl-d14	1.000	1.022	-2.2	55	0.00
30		Benzo(a)anthracene	1.000	0.904	9.6	48	0.00
31		3,3'-Dichlorobenzidine	1.000	0.810	19.0	45	0.00
32		Chrysene	1.000	0.985	1.5	53	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.861	13.9	42	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.896	10.4	59	0.01
35	I	Perylene-d12	5.000	5.000	0.0	52	0.00
36		Benzo(b)fluoranthene	1.000	0.928	7.2	51	0.00
37		Benzo(k)fluoranthene	1.000	0.918	8.2	50	0.00
38	C	Benzo(a)pyrene	1.000	0.883	11.7	50	0.00
39		Dibenzo(a,h)anthracene	1.000	0.928	7.2	61	0.00
40		Benzo(g,h,i)perylene	1.000	0.914	8.6	61	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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