Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040915\
Data File : BM000935.D
Acq On : 09 Apr 2015 23:09
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040915

Quant Time: Apr 10 16:34:28 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 10 16:15:17 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	103	0.00
2		1,4-Dioxane	0.497	0.499	-0.4	106	0.00
3		n-Nitrosodimethylamine	0.486	0.498	-2.5	108	0.00
4	S	2-Fluorophenol	1.058	1.125	-6.3	109	0.00
5	S	Phenol-d6	1.267	1.343	-6.0	112	0.00
6	C	Phenol	1.302	1.367	-5.0	111	0.00
7		bis(2-Chloroethyl)ether	1.142	1.215	-6.4	110	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
9	S	Nitrobenzene-d5	0.241	0.259	-7.5	112	0.00
10		Nitrobenzene	0.283	0.309	-9.2	112	0.00
11		2,4-Dimethylphenol	0.275	0.295	-7.3	110	0.00
12	C	2,4-Dichlorophenol	0.243	0.257	-5.8	112	0.00
13		Naphthalene	1.009	1.058	-4.9	107	0.00
14	C	Hexachlorobutadiene	0.184	0.197	-7.1	109	0.00
15	C	2-Methylnaphthalene	0.689	0.738	-7.1	109	0.00
16		1-Methylnaphthalene	0.694	0.751	-8.2	110	-0.01
17	I	Acenaphthene-d10	1.000	1.000	0.0	125	0.01
18	S	2,4,6-Tribromophenol	0.179	0.164	8.4	93	0.01
19	S	2-Fluorobiphenyl	1.401	1.437	-2.6	110	0.00
20		Acenaphthylene	2.032	1.913	5.9	88	0.01
21	C	Acenaphthene	1.380	1.460	-5.8	123	-0.02
22	P	2,4-Dinitrophenol	0.058	0.046#	20.7	140	-0.01
23		Fluorene	1.606	1.810	-12.7	134	-0.02
24	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.01
25		Hexachlorobenzene	0.270	0.277	-2.6	105	0.01
26	C	Pentachlorophenol	0.086	0.080	7.0	111	0.00
27		Phenanthrene	1.292	1.325	-2.6	104	0.01
28		Anthracene	1.184	1.248	-5.4	117	-0.01
29	C	Fluoranthene	1.350	1.409	-4.4	108	-0.01
30	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
31		Benzidine	0.300	0.295	1.7	121	0.00
32		Pyrene	1.625	1.737	-6.9	108	0.00
33	S	Terphenyl-d14	0.609	0.669	-9.9	114	0.00
34		Benzo(a)anthracene	1.443	1.545	-7.1	113	-0.01
35		3,3'-Dichlorobenzidine	0.389	0.412	-5.9	119	0.00
36		Chrysene	1.475	1.555	-5.4	106	0.00
37		Indeno(1,2,3-cd)pyrene	1.506	1.642	-9.0	112	0.00
38	I	Perylene-d12	1.000	1.000	0.0	102	0.00
39		Benzo(b)fluoranthene	1.565	1.754	-12.1	120	0.00
40		Benzo(k)fluoranthene	1.506	1.534	-1.9	101	0.00
41	C	Benzo(a)pyrene	1.378	1.468	-6.5	111	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Dibenzo(a,h)anthracene	1.278	1.412	-10.5	114	0.00
43	Benzo(g,h,i)perylene	1.410	1.514	-7.4	110	0.00

(#) = Out of Range SPCC's out = 1 CCC's out = 0