Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004483.D
Acq On : 29 Feb 2016 22:24
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM022916

Quant Time: Mar 01 05:58:08 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 01 00:52:28 2016
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	96	0.00
2		1,4-Dioxane	0.580	0.483	16.7	94	0.00
3		n-Nitrosodimethylamine	0.508	0.479	5.7	96	0.00
4	S	2-Fluorophenol	1.162	1.119	3.7	98	0.00
5	S	Phenol-d6	1.313	1.299	1.1	100	0.00
6		bis(2-Chloroethyl)ether	1.389	1.394	-0.4	99	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
8	S	Nitrobenzene-d5	0.254	0.246	3.1	101	0.00
9		Nitrobenzene	0.344	0.331	3.8	99	0.00
10		Naphthalene	1.572	1.392	11.5	100	0.00
11		Hexachlorobutadiene	0.230	0.228	0.9	100	-0.01
12		2-Methylnaphthalene	0.869	0.868	0.1	103	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
14	S	2,4,6-Tribromophenol	0.192	0.182	5.2	104	0.00
15	S	2-Fluorobiphenyl	1.531	1.553	-1.4	105	0.00
16		Acenaphthylene	2.861	2.749	3.9	99	0.00
17		Acenaphthene	1.872	1.897	-1.3	111	0.00
18		Fluorene	2.285	2.381	-4.2	109	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
20		4-Bromophenyl-phenylether	0.274	0.294	-7.3	122	0.00
21		Hexachlorobenzene	0.308	0.356	-15.6	129	0.00
22		Pentachlorophenol	0.074	0.058	21.6	120	0.00
23		Phenanthrene	1.762	1.926	-9.3	121	0.00
24		Anthracene	1.485	1.479	0.4	99	0.02
25		Fluoranthene	1.891	1.993	-5.4	107	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	87	0.01
27		Benzidine	0.535	0.995	-86.0#	206#	0.00
28		Pyrene	1.977	2.123	-7.4	107	0.00
29	S	Terphenyl-d14	0.643	0.685	-6.5	106	0.00
30		Benzo(a)anthracene	1.845	1.722	6.7	91	0.01
31		3,3'-Dichlorobenzidine	1.118	1.660	-48.5#	163#	0.00
32		Chrysene	3.757	2.233	40.6#	90	0.00
33		Bis(2-ethylhexyl)phthalate	1.432	1.170	18.3	68	-0.01
34		Indeno(1,2,3-cd)pyrene	1.712	1.785	-4.3	101	-0.01
35	I	Perylene-d12	1.000	1.000	0.0	99	-0.01
36		Benzo(b)fluoranthene	1.949	1.922	1.4	98	0.00
37		Benzo(k)fluoranthene	1.920	1.798	6.4	97	0.00
38	C	Benzo(a)pyrene	1.761	1.772	-0.6	102	0.00
39		Dibenzo(a,h)anthracene	1.533	1.497	2.3	100	0.00
40		Benzo(g,h,i)perylene	1.683	1.659	1.4	101	-0.01

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004483.D
Acq On : 29 Feb 2016 22:24
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM022916

Quant Time: Mar 01 05:58:08 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 01 00:52:28 2016
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)

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