Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022816\
Data File : BM004460.D
Acq On : 28 Feb 2016 20:56
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM022816

Quant Time: Feb 29 06:36:15 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM022816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 29 05:57:18 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	86	0.00
2		1,4-Dioxane	0.475	0.453	4.6	104	0.00
3		n-Nitrosodimethylamine	0.371	0.376	-1.3	101	0.00
4	S	2-Fluorophenol	1.100	1.087	1.2	97	0.00
5	S	Phenol-d6	1.248	1.141	8.6	97	0.00
6		bis(2-Chloroethyl)ether	0.992	1.024	-3.2	100	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
8	S	Nitrobenzene-d5	0.276	0.258	6.5	97	0.00
9		Nitrobenzene	0.307	0.296	3.6	98	0.00
10		Hexachlorobutadiene	0.204	0.201	1.5	98	0.00
11		2-Methylnaphthalene	0.761	0.778	-2.2	103	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	0.198	0.150	24.2	89	0.00
14	S	2-Fluorobiphenyl	1.525	1.430	6.2	99	0.00
15	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
16		Hexachlorobenzene	0.278	0.268	3.6	104	0.00
17		Pentachlorophenol	0.103	0.065	36.9#	94	0.00
18		Fluoranthene	1.981	1.765	10.9	100	0.00
19	I	Chrysene-d12	1.000	1.000	0.0	93	0.00
20		Benzidine	0.257	0.361	-40.5#	158#	0.00
21		Pyrene	1.489	1.400	6.0	99	0.00
22	S	Terphenyl-d14	0.637	0.616	3.3	103	0.00
23		Benzo(a)anthracene	1.428	1.268	11.2	105	0.00
24		3,3'-Dichlorobenzidine	0.490	0.512	-4.5	101	0.00
25		Chrysene	2.306	1.355	41.2#	99	0.00
26		Bis(2-ethylhexyl)phthalate	0.906	0.696	23.2	86	0.00
27		Indeno(1,2,3-cd)pyrene	1.301	1.234	5.1	102	-0.01
28	I	Perylene-d12	1.000	1.000	0.0	91	-0.01
29		Benzo(b)fluoranthene	1.885	1.794	4.8	96	0.00
30		Benzo(k)fluoranthene	1.625	1.448	10.9	100	0.00
31	C	Benzo(a)pyrene	1.466	1.453	0.9	102	0.00
32		Dibenzo(a,h)anthracene	1.309	1.250	4.5	102	0.00
33		Benzo(g,h,i)perylene	1.438	1.380	4.0	103	-0.01

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