Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM120715\
Data File : BM003407.D
Acq On : 07 Dec 2015 14:48
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM120715

Quant Time: Dec 07 15:30:40 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM120715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 07 15:08:30 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	114	0.00
2		1,4-Dioxane	0.449	0.391	12.9	109	0.00
3		n-Nitrosodimethylamine	0.394	0.385	2.3	114	0.00
4	S	2-Fluorophenol	0.857	0.832	2.9	118	-0.02
5	S	Phenol-d6	1.203	1.204	-0.1	120	0.00
6		2-Chlorophenol	1.120	1.127	-0.6	119	0.00
7	C	Phenol	1.258	1.240	1.4	118	0.00
8		bis(2-Chloroethyl)ether	0.896	0.903	-0.8	118	0.00
9		2-Methylphenol	0.708	0.710	-0.3	121	0.00
10		3+4-Methylphenols	1.029	1.034	-0.5	119	0.00
11	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
12	S	Nitrobenzene-d5	0.264	0.270	-2.3	122	0.00
13		Nitrobenzene	0.275	0.279	-1.5	123	0.00
14	C	2-Nitrophenol	0.112	0.110	1.8	124	0.00
15		2,4-Dimethylphenol	0.264	0.269	-1.9	122	0.00
16	C	2,4-Dichlorophenol	0.246	0.251	-2.0	119	0.00
17		Hexachlorobutadiene	0.183	0.175	4.4	113	0.00
18	C	4-Chloro-3-methylphenol	0.214	0.211	1.4	119	0.03
19		2-Methylnaphthalene	0.773	0.764	1.2	118	0.00
20	I	Acenaphthene-d10	1.000	1.000	0.0	145	0.00
21	S	2,4,6-Tribromophenol	0.123	0.096	22.0	101	0.00
22	C	2,4,6-Trichlorophenol	0.335	0.280	16.4	127	0.00
23		2,4,5-Trichlorophenol	0.368	0.308	16.3	116	0.00
24	S	2-Fluorobiphenyl	1.760	1.534	12.8	124	0.00
25	P	2,4-Dinitrophenol	0.038	0.037#	2.6	137	0.00
26	P	4-Nitrophenol	0.153	0.141	7.8	120	0.00
27	I	Phenanthrene-d10	1.000	1.000	0.0	133	0.00
28		4,6-Dinitro-2-methylphenol	0.061	0.064	-4.9	147	0.00
29		Hexachlorobenzene	0.212	0.182	14.2	107	0.00
30	C	Pentachlorophenol	0.065	0.049	24.6#	120	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	85	0.00
32		Benzidine	0.323	0.379	-17.3	126	0.00
33	S	Terphenyl-d14	0.676	0.794	-17.5	104	0.00
34		3,3'-Dichlorobenzidine	0.304	0.348	-14.5	134	0.00
35	I	Perylene-d12	1.000	1.000	0.0	114	0.01

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