Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM113015\
Data File : BM003335.D
Acq On : 01 Dec 2015 00:57
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 01 02:52:46 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM112715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 27 16:36:40 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	134	-0.02
2		1,4-Dioxane	0.488	0.457	6.4	130	-0.02
3		n-Nitrosodimethylamine	0.462	0.492	-6.5	143	-0.02
4	S	2-Fluorophenol	0.986	0.991	-0.5	139	0.00
5	S	Phenol-d6	1.131	1.229	-8.7	151#	-0.02
6		bis(2-Chloroethyl)ether	1.156	1.246	-7.8	148	-0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	145	-0.02
8	S	Nitrobenzene-d5	0.222	0.238	-7.2	164#	-0.01
9		Nitrobenzene	0.243	0.260	-7.0	165#	-0.01
10		Naphthalene	1.081	1.092	-1.0	149	0.00
11		Hexachlorobutadiene	0.164	0.174	-6.1	158#	-0.02
12		2-Methylnaphthalene	0.678	0.744	-9.7	162#	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	166#	0.00
14	S	2,4,6-Tribromophenol	0.090	0.111	-23.3	210#	0.00
15	S	2-Fluorobiphenyl	1.614	1.500	7.1	167#	0.00
16		Acenaphthylene	1.832	2.010	-9.7	190#	0.00
17		Acenaphthene	1.414	1.292	8.6	176#	0.00
18		Fluorene	1.427	1.455	-2.0	177#	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	162#	0.00
20		4-Bromophenyl-phenylether	0.170	0.139	18.2	179#	-0.03
21		Hexachlorobenzene	0.176	0.178	-1.1	180#	0.00
22		Pentachlorophenol	0.066	0.053	19.7	189#	0.00
23		Phenanthrene	1.145	0.909	20.6	161#	0.00
24		Anthracene	0.933	1.034	-10.8	179#	0.00
25		Fluoranthene	1.090	0.910	16.5	151#	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	114	-0.01
27		Benzidine	0.229	0.394	-72.1#	275#	0.00
28		Pyrene	1.528	1.706	-11.6	151#	0.00
29	S	Terphenyl-d14	0.697	0.802	-15.1	144	0.00
30		Benzo(a)anthracene	1.266	1.470	-16.1	149	-0.01
31		3,3'-Dichlorobenzidine	0.270	0.295	-9.3	183#	0.00
32		Chrysene	1.515	1.281	15.4	116	0.00
33		Bis(2-ethylhexyl)phthalate	0.518	0.532	-2.7	193#	0.00
34		Indeno(1,2,3-cd)pyrene	1.160	1.175	-1.3	133	-0.01
35	I	Perylene-d12	1.000	1.000	0.0	116	-0.01
36		Benzo(b)fluoranthene	1.439	1.468	-2.0	116	0.00
37		Benzo(k)fluoranthene	1.294	1.369	-5.8	129	-0.01
38	C	Benzo(a)pyrene	1.146	1.197	-4.5	127	-0.01
39		Dibenzo(a,h)anthracene	0.992	1.192	-20.2	141	-0.01
40		Benzo(g,h,i)perylene	1.190	1.246	-4.7	125	-0.01

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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