Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE070516\
Data File : BE091967.D
Acq On : 5 Jul 2016 15:55
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE070516

Quant Time: Jul 05 16:51:49 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE070516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 05 16:14:58 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	91	0.00
2		1,4-Dioxane	0.678	0.736	-8.6	95	0.00
3		n-Nitrosodimethylamine	0.768	0.731	4.8	93	-0.02
4	S	2-Fluorophenol	1.196	1.139	4.8	92	0.00
5	S	Phenol-d6	1.662	1.577	5.1	95	0.00
6		bis(2-Chloroethyl)ether	1.449	1.480	-2.1	101	-0.02
7		n-Nitroso-di-n-propylamine	1.387	1.327	4.3	94	-0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
9	S	Nitrobenzene-d5	0.341	0.330	3.2	91	-0.02
10		Nitrobenzene	0.370	0.348	5.9	93	0.00
11		bis(2-Chloroethoxy)methane	0.418	0.418	0.0	81	0.00
12		Naphthalene	1.164	1.201	-3.2	93	0.00
13		Hexachlorobutadiene	0.177	0.185	-4.5	91	0.00
14		2-Methylnaphthalene	0.748	0.775	-3.6	94	-0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
16	S	2,4,6-Tribromophenol	0.143	0.128	10.5	94	0.00
17	S	2-Fluorobiphenyl	1.220	1.340	-9.8	95	0.00
18		Acenaphthylene	6.276	6.350	-1.2	93	0.00
19		2,6-Dinitrotoluene	1.093	0.857	21.6	96	0.00
20		Acenaphthene	1.367	1.497	-9.5	95	0.00
21		2,4-Dinitrotoluene	1.115	0.833	25.3#	101	0.00
22		Fluorene	1.442	1.571	-8.9	94	0.00
23		4-Chlorophenyl-phenylether	0.719	0.817	-13.6	96	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
25		4-Bromophenyl-phenylether	0.217	0.223	-2.8	96	0.00
26		Hexachlorobenzene	0.216	0.227	-5.1	98	0.00
27		Pentachlorophenol	0.078	0.047	39.7#	100	0.00
28		Phenanthrene	1.303	1.379	-5.8	97	0.00
29		Anthracene	1.218	1.238	-1.6	97	0.00
30		Fluoranthene	5.599	5.958	-6.4	102	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	97	0.00
32		Benzidine	0.333	0.262	21.3	90	0.00
33		Pyrene	3.640	3.750	-3.0	99	0.00
34	S	Terphenyl-d14	0.590	0.653	-10.7	106	0.00
35		Benzo(a)anthracene	0.837	0.863	-3.1	101	0.00
36		3,3'-Dichlorobenzidine	0.257	0.249	3.1	94	0.00
37		Chrysene	1.187	1.263	-6.4	104	0.00
38		Bis(2-ethylhexyl)phthalate	0.970	0.896	7.6	98	0.00
39		Indeno(1,2,3-cd)pyrene	4.657	5.008	-7.5	101	0.00
40	I	Perylene-d12	1.000	1.000	0.0	97	0.00
41		Benzo(b)fluoranthene	1.381	1.388	-0.5	99	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.335	1.419	-6.3	105	0.00
43 C	Benzo(a)pyrene	1.304	1.366	-4.8	103	0.00
44	Dibenzo(a,h)anthracene	1.227	1.254	-2.2	102	0.00
45	Benzo(g,h,i)perylene	5.004	5.096	-1.8	100	0.02

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