Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE061416\
Data File : BE091917.D
Acq On : 14 Jun 2016 19:37
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE061416

Quant Time: Jun 15 12:18:22 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE061416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 15 11:58:50 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	105	0.00
2		1,4-Dioxane	0.705	0.735	-4.3	106	0.00
3		n-Nitrosodimethylamine	0.802	0.808	-0.7	105	0.02
4	S	2-Fluorophenol	0.977	0.903	7.6	98	0.00
5	S	Phenol-d6	1.267	1.156	8.8	100	0.02
6		bis(2-Chloroethyl)ether	1.431	1.451	-1.4	107	0.02
7		n-Nitroso-di-n-propylamine	1.137	1.100	3.3	103	0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
9	S	Nitrobenzene-d5	0.345	0.340	1.4	106	0.00
10		Nitrobenzene	0.343	0.335	2.3	104	0.02
11		bis(2-Chloroethoxy)methane	0.438	0.475	-8.4	109	0.00
12		Naphthalene	1.092	1.176	-7.7	109	0.00
13		Hexachlorobutadiene	0.177	0.187	-5.6	107	0.00
14		2-Methylnaphthalene	0.741	0.759	-2.4	107	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
16	S	2,4,6-Tribromophenol	0.102	0.089	12.7	101	0.00
17	S	2-Fluorobiphenyl	1.269	1.339	-5.5	107	0.00
18		Acenaphthylene	5.075	4.718	7.0	102	0.03
19		2,6-Dinitrotoluene	0.811	0.614	24.3	117	0.00
20		Acenaphthene	1.398	1.427	-2.1	111	0.00
21		2,4-Dinitrotoluene	0.758	0.632	16.6	97	0.03
22		Fluorene	1.391	1.390	0.1	108	0.02
23		4-Chlorophenyl-phenylether	0.685	0.710	-3.6	107	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
25		4-Bromophenyl-phenylether	0.193	0.203	-5.2	110	0.02
26		Hexachlorobenzene	0.198	0.209	-5.6	108	0.00
27		Pentachlorophenol	0.054	0.040	25.9#	99	0.02
28		Phenanthrene	1.250	1.332	-6.6	109	0.00
29		Anthracene	1.106	1.115	-0.8	109	0.00
30		Fluoranthene	5.354	5.468	-2.1	107	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
32		Benzidine	0.170	0.113	33.5#	88	0.00
33		Pyrene	3.401	3.884	-14.2	111	0.00
34	S	Terphenyl-d14	0.612	0.732	-19.6	110	0.00
35		Benzo(a)anthracene	0.786	0.757	3.7	93	0.00
36		3,3'-Dichlorobenzidine	0.127	0.107	15.7	95	0.03
37		Chrysene	1.147	1.184	-3.2	88	0.00
38		Bis(2-ethylhexyl)phthalate	0.580	0.554	4.5	104	0.00
39		Indeno(1,2,3-cd)pyrene	3.925	4.104	-4.6	100	0.00
40	I	Perylene-d12	1.000	1.000	0.0	100	0.00
41		Benzo(b)fluoranthene	1.284	1.313	-2.3	98	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.175	1.213	-3.2	103	0.00
43 C	Benzo(a)pyrene	1.122	1.092	2.7	98	0.00
44	Dibenzo(a,h)anthracene	1.092	1.107	-1.4	101	0.00
45	Benzo(g,h,i)perylene	4.507	4.488	0.4	98	0.00

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