Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE061516\
Data File : BE091933.D
Acq On : 15 Jun 2016 20:14
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 16 01:00:28 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 12:00:38 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	82	0.02
2		1,4-Dioxane	0.705	0.706	-0.1	80	0.00
3		n-Nitrosodimethylamine	0.802	0.632	21.2	64	0.04
4	S	2-Fluorophenol	0.977	0.853	12.7	72	0.02
5	S	Phenol-d6	1.267	1.092	13.8	74	0.02
6		bis(2-Chloroethyl)ether	1.431	1.232	13.9	71	0.02
7		n-Nitroso-di-n-propylamine	1.137	1.002	11.9	73	0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	84	0.03
9	S	Nitrobenzene-d5	0.345	0.314	9.0	76	0.02
10		Nitrobenzene	0.343	0.300	12.5	72	0.02
11		bis(2-Chloroethoxy)methane	0.438	0.372	15.1	66	0.03
12		Naphthalene	1.092	1.055	3.4	76	0.00
13		Hexachlorobutadiene	0.177	0.170	4.0	75	0.00
14		2-Methylnaphthalene	0.741	0.682	8.0	75	0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	79	0.00
16	S	2,4,6-Tribromophenol	0.102	0.082	19.6	68	0.02
17	S	2-Fluorobiphenyl	1.269	1.276	-0.6	74	0.00
18		Acenaphthylene	5.075	4.541	10.5	71	0.03
19		2,6-Dinitrotoluene	0.811	0.503	38.0#	69	0.00
20		Acenaphthene	1.398	1.305	6.7	74	0.00
21		2,4-Dinitrotoluene	0.758	0.667	12.0	74	0.03
22		Fluorene	1.391	1.270	8.7	72	0.02
23		4-Chlorophenyl-phenylether	0.685	0.636	7.2	70	0.02
24	I	Phenanthrene-d10	1.000	1.000	0.0	78	0.00
25		4-Bromophenyl-phenylether	0.193	0.177	8.3	70	0.02
26		Hexachlorobenzene	0.198	0.194	2.0	73	0.00
27		Pentachlorophenol	0.054	0.030	44.4#	55	0.02
28		Phenanthrene	1.250	1.206	3.5	72	0.00
29		Anthracene	1.106	0.990	10.5	71	0.00
30		Fluoranthene	5.354	4.479	16.3	64	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
32		Benzidine	0.170	0.107	37.1#	70	0.02
33		Pyrene	3.401	2.831	16.8	67	0.00
34	S	Terphenyl-d14	0.612	0.488	20.3	61	0.00
35		Benzo(a)anthracene	0.786	0.775	1.4	80	0.00
36		3,3'-Dichlorobenzidine	0.127	0.090	29.1#	67	0.03
37		Chrysene	1.147	1.185	-3.3	74	0.00
38		Bis(2-ethylhexyl)phthalate	0.580	0.453	21.9	71	-0.03
39		Indeno(1,2,3-cd)pyrene	3.925	3.209	18.2	66	0.00
40	I	Perylene-d12	1.000	1.000	0.0	75	-0.01
41		Benzo(b)fluoranthene	1.284	1.225	4.6	69	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.175	1.001	14.8	64	0.00
43 C	Benzo(a)pyrene	1.122	0.974	13.2	66	0.00
44	Dibenzo(a,h)anthracene	1.092	0.952	12.8	65	-0.02
45	Benzo(g,h,i)perylene	4.507	3.952	12.3	65	-0.02

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