Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE122617\
Data File : BE095047.D
Acq On : 26 Dec 2017 9:39
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 26 10:52:14 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 14 18:42:53 2017
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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	86	-0.01
2		1,4-Dioxane	0.400	0.383	4.3	77	0.00
3		n-Nitrosodimethylamine	0.400	0.405	-1.3	86	0.00
4	S	2-Fluorophenol	0.400	0.389	2.8	87	-0.01
5	S	Phenol-d6	0.400	0.380	5.0	84	0.00
6		bis(2-Chloroethyl)ether	0.400	0.412	-3.0	89	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	93	-0.02
8	S	Nitrobenzene-d5	0.400	0.379	5.3	92	0.00
9		Naphthalene	0.400	0.381	4.8	89	-0.02
10		Hexachlorobutadiene	0.400	0.385	3.8	89	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.384	4.0	90	-0.02
12		2-Methylnaphthalene	0.400	0.348	13.0	82	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	89	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.323	19.3	77	-0.03
15	S	2-Fluorobiphenyl	0.400	0.394	1.5	87	-0.02
16		Acenaphthylene	0.400	0.371	7.3	87	-0.03
17		Acenaphthene	0.400	0.385	3.8	87	-0.02
18		Fluorene	0.400	0.384	4.0	88	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	90	-0.03
20		4-Bromophenyl-phenylether	0.400	0.374	6.5	87	-0.03
21		Hexachlorobenzene	0.400	0.384	4.0	88	-0.03
22		Pentachlorophenol	0.400	0.348	13.0	81	-0.02
23		Phenanthrene	0.400	0.388	3.0	90	-0.03
24		Anthracene	0.400	0.374	6.5	89	-0.03
25	SURR	Fluoranthene-d10	0.400	0.387	3.3	91	-0.03
26		Fluoranthene	0.400	0.387	3.3	90	-0.03
27	I	Chrysene-d12	0.400	0.400	0.0	89	-0.03
28		Pyrene	0.400	0.489	-22.2	105	-0.03
29	S	Terphenyl-d14	0.400	0.400	0.0	88	-0.03
30		Benzo(a)anthracene	0.400	0.408	-2.0	92	-0.03
31		Chrysene	0.400	0.418	-4.5	93	-0.03
32		Bis(2-ethylhexyl)phthalate	0.400	0.356	11.0	89	-0.05
33		Indeno(1,2,3-cd)pyrene	0.400	0.423	-5.7	92	-0.07
34	I	Perylene-d12	0.400	0.400	0.0	98	-0.05
35		Benzo(b)fluoranthene	0.400	0.388	3.0	95	-0.04
36		Benzo(k)fluoranthene	0.400	0.378	5.5	97	-0.05
37	C	Benzo(a)pyrene	0.400	0.395	1.3	97	-0.05
38		Dibenzo(a,h)anthracene	0.400	0.343	14.2	85	-0.07
39		Benzo(g,h,i)perylene	0.400	0.375	6.3	94	-0.08

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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