Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101717\
Data File : BE094435.D
Acq On : 17 Oct 2017 18:39
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 17 20:35:49 2017
Quant Method : Z:\HPCHEM1\BNA_E\Methods\8270-SIM-BE101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 14:54:40 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	96	-0.02
2		1,4-Dioxane	0.400	0.426	-6.5	98	0.00
3		n-Nitrosodimethylamine	0.400	0.345	13.8	83	0.00
4	S	2-Fluorophenol	0.400	0.387	3.3	91	0.00
5	S	Phenol-d6	0.400	0.383	4.3	89	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.393	1.8	91	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	95	0.00
8	S	Nitrobenzene-d5	0.400	0.384	4.0	92	0.00
9		Naphthalene	0.400	0.399	0.3	95	-0.02
10		Hexachlorobutadiene	0.400	0.380	5.0	91	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	97	0.00
12		2-Methylnaphthalene	0.400	0.410	-2.5	99	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14	S	2,4,6-Tribromophenol	0.400	0.363	9.3	90	0.00
15	S	2-Fluorobiphenyl	0.400	0.385	3.8	99	0.00
16		Acenaphthylene	0.400	0.393	1.8	103	0.00
17		Acenaphthene	0.400	0.390	2.5	103	0.00
18		Fluorene	0.400	0.397	0.8	105	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	98	-0.03
20		4-Bromophenyl-phenylether	0.400	0.418	-4.5	114	-0.03
21		Hexachlorobenzene	0.400	0.356	11.0	96	0.00
22		Pentachlorophenol	0.400	0.379	5.3	78	0.03
23		Phenanthrene	0.400	0.349	12.8	82	0.00
24		Anthracene	0.400	0.408	-2.0	103	0.00
25	SURR	Fluoranthene-d10	0.400	0.398	0.5	101	0.00
26		Fluoranthene	0.400	0.400	0.0	103	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	109	0.00
28		Pyrene	0.400	0.369	7.8	102	0.00
29	S	Terphenyl-d14	0.400	0.374	6.5	101	0.00
30		Benzo(a)anthracene	0.400	0.386	3.5	106	-0.01
31		Chrysene	0.400	0.381	4.8	102	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.359	10.3	102	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.384	4.0	105	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	109	-0.02
35		Benzo(b)fluoranthene	0.400	0.379	5.3	103	0.00
36		Benzo(k)fluoranthene	0.400	0.404	-1.0	109	0.00
37	C	Benzo(a)pyrene	0.400	0.387	3.3	105	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.384	4.0	103	-0.01
39		Benzo(g,h,i)perylene	0.400	0.389	2.8	105	-0.02

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

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