Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\BE040418\
Data File : BE095939.D
Acq On : 4 Apr 2018 10:06
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 05 02:46:24 2018
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 23 12:41:11 2018
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	122	-0.01
2		1,4-Dioxane	0.400	0.362	9.5	116	0.00
3		n-Nitrosodimethylamine	0.400	0.384	4.0	113	0.00
4	S	2-Fluorophenol	0.400	0.375	6.3	110	-0.01
5	S	Phenol-d6	0.400	0.423	-5.7	123	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.413	-3.2	121	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	129	0.00
8	S	Nitrobenzene-d5	0.400	0.393	1.8	125	0.00
9		Naphthalene	0.400	0.420	-5.0	130	-0.01
10		Hexachlorobutadiene	0.400	0.375	6.3	115	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	129	-0.01
12		2-Methylnaphthalene	0.400	0.423	-5.7	132	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	127	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.354	11.5	116	-0.01
15	S	2-Fluorobiphenyl	0.400	0.406	-1.5	126	-0.01
16		Acenaphthylene	0.400	0.418	-4.5	133	0.00
17		Acenaphthene	0.400	0.413	-3.2	132	0.00
18		Fluorene	0.400	0.409	-2.2	128	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	120	-0.01
20		4-Bromophenyl-phenylether	0.400	0.404	-1.0	120	0.00
21		Hexachlorobenzene	0.400	0.410	-2.5	121	0.00
22		Pentachlorophenol	0.400	0.473	-18.2	162	-0.01
23		Phenanthrene	0.400	0.418	-4.5	122	0.00
24		Anthracene	0.400	0.406	-1.5	121	-0.01
25	SURR	Fluoranthene-d10	0.400	0.374	6.5	110	0.00
26		Fluoranthene	0.400	0.382	4.5	114	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	109	-0.01
28		Pyrene	0.400	0.463	-15.8	114	-0.01
29	S	Terphenyl-d14	0.400	0.401	-0.3	108	0.00
30		Benzo(a)anthracene	0.400	0.409	-2.2	109	0.00
31		Chrysene	0.400	0.418	-4.5	112	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.367	8.3	106	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.415	-3.7	111	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	110	-0.02
35		Benzo(b)fluoranthene	0.400	0.413	-3.2	112	-0.01
36		Benzo(k)fluoranthene	0.400	0.419	-4.7	114	-0.01
37	C	Benzo(a)pyrene	0.400	0.413	-3.2	111	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.411	-2.7	110	-0.02
39		Benzo(g,h,i)perylene	0.400	0.417	-4.2	112	-0.02

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

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