Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE062217\
Data File : BE093332.D
Acq On : 22 Jun 2017 17:25
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 23 08:10:53 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 12 14:07:54 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	28	-0.02
2		1,4-Dioxane	0.400	0.352	12.0	30	-0.01
3		n-Nitrosodimethylamine	0.400	0.389	2.8	30	-0.01
4	S	2-Fluorophenol	0.400	0.451	-12.7	37	0.00
5	S	Phenol-d6	0.400	0.503	-25.7#	41	0.00
6		bis(2-Chloroethyl)ether	0.400	0.410	-2.5	33	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	32	-0.02
8	S	Nitrobenzene-d5	0.400	0.460	-15.0	43	-0.02
9		Naphthalene	0.400	0.420	-5.0	38	0.00
10		Hexachlorobutadiene	0.400	0.412	-3.0	38	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.465	-16.3	43	-0.01
12		2-Methylnaphthalene	0.400	0.457	-14.2	42	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	39	0.00
14	S	2,4,6-Tribromophenol	0.400	0.534	-33.5#	62	0.00
15	S	2-Fluorobiphenyl	0.400	0.402	-0.5	43	0.00
16		Acenaphthylene	0.400	0.452	-13.0	52	0.00
17		Acenaphthene	0.400	0.423	-5.7	47	-0.02
18		Fluorene	0.400	0.431	-7.7	48	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	35	-0.03
20		4-Bromophenyl-phenylether	0.400	0.299	25.3#	30	0.00
21		Hexachlorobenzene	0.400	0.518	-29.5#	58	0.00
22		Pentachlorophenol	0.400	0.474	-18.5	52	0.00
23		Phenanthrene	0.400	0.325	18.8	33	0.00
24		Anthracene	0.400	0.398	0.5	45	0.00
25	SURR	Fluoranthene-d10	0.400	0.413	-3.2	43	0.00
26		Fluoranthene	0.400	0.390	2.5	42	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	28	-0.02
28		Pyrene	0.400	0.433	-8.2	33	0.00
29	S	Terphenyl-d14	0.400	0.437	-9.2	34	0.00
30		Benzo(a)anthracene	0.400	0.473	-18.2	37	0.00
31		Chrysene	0.400	0.429	-7.2	33	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.497	-24.2	37	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.425	-6.2	34	0.00
34	I	Perylene-d12	0.400	0.400	0.0	26	-0.01
35		Benzo(b)fluoranthene	0.400	0.422	-5.5	32	-0.01
36		Benzo(k)fluoranthene	0.400	0.453	-13.2	34	-0.01
37	C	Benzo(a)pyrene	0.400	0.444	-11.0	34	0.00
38		Dibenzo(a,h)anthracene	0.400	0.429	-7.2	33	0.00
39		Benzo(g,h,i)perylene	0.400	0.423	-5.7	32	0.02

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

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