Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120817\
Data File : BE094969.D
Acq On : 9 Dec 2017 3:21
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 09 05:46:40 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 01 13:33:28 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	84	0.00
2		1,4-Dioxane	0.400	0.373	6.8	74	0.00
3		n-Nitrosodimethylamine	0.400	0.357	10.8	76	-0.01
4	S	2-Fluorophenol	0.400	0.397	0.8	83	0.00
5	S	Phenol-d6	0.400	0.401	-0.3	84	0.00
6		bis(2-Chloroethyl)ether	0.400	0.367	8.3	75	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	81	0.00
8	S	Nitrobenzene-d5	0.400	0.559	-39.8#	114	0.00
9		Naphthalene	0.400	0.397	0.8	80	0.00
10		Hexachlorobutadiene	0.400	0.405	-1.3	81	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	82	0.00
12		2-Methylnaphthalene	0.400	0.400	0.0	82	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	84	0.00
14	S	2,4,6-Tribromophenol	0.400	0.518	-29.5#	115	0.00
15	S	2-Fluorobiphenyl	0.400	0.400	0.0	83	0.00
16		Acenaphthylene	0.400	0.372	7.0	81	0.00
17		Acenaphthene	0.400	0.391	2.3	84	0.00
18		Fluorene	0.400	0.391	2.3	81	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	84	0.00
20		4-Bromophenyl-phenylether	0.400	0.382	4.5	81	0.00
21		Hexachlorobenzene	0.400	0.412	-3.0	85	0.01
22		Pentachlorophenol	0.400	0.437	-9.2	94	0.00
23		Phenanthrene	0.400	0.384	4.0	80	0.00
24		Anthracene	0.400	0.363	9.3	78	0.00
25	SURR	Fluoranthene-d10	0.400	0.356	11.0	75	0.00
26		Fluoranthene	0.400	0.357	10.8	75	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	68	0.00
28		Pyrene	0.400	0.429	-7.2	74	0.00
29	S	Terphenyl-d14	0.400	0.432	-8.0	74	0.00
30		Benzo(a)anthracene	0.400	0.401	-0.3	70	-0.01
31		Chrysene	0.400	0.372	7.0	63	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.368	8.0	73	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.413	-3.2	71	0.00
34	I	Perylene-d12	0.400	0.400	0.0	71	0.00
35		Benzo(b)fluoranthene	0.400	0.388	3.0	70	0.00
36		Benzo(k)fluoranthene	0.400	0.379	5.3	69	0.00
37	C	Benzo(a)pyrene	0.400	0.392	2.0	72	0.00
38		Dibenzo(a,h)anthracene	0.400	0.394	1.5	72	-0.01
39		Benzo(g,h,i)perylene	0.400	0.389	2.8	70	-0.01

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

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