Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070519\
Data File : BE100375.D
Acq On : 5 Jul 2019 11:18
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 06 00:10:31 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 14:06:53 2019
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	76	-0.02
2		1,4-Dioxane	0.400	0.430	-7.5	79	-0.02
3		n-Nitrosodimethylamine	0.400	0.563	-40.7#	129	-0.04
4	S	2-Fluorophenol	0.400	0.479	-19.7	90	-0.02
5	S	Phenol-d6	0.400	0.343	14.2	65	0.00
6		bis(2-Chloroethyl)ether	0.400	0.351	12.3	60	-0.03
7	I	Naphthalene-d8	0.400	0.400	0.0	67	-0.03
8	S	Nitrobenzene-d5	0.400	0.322	19.5	55	-0.01
9		Naphthalene	0.400	0.400	0.0	66	-0.01
10		Hexachlorobutadiene	0.400	0.471	-17.7	76	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.390	2.5	64	-0.01
12		2-Methylnaphthalene	0.400	0.350	12.5	62	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	71	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.382	4.5	69	0.00
15	S	2-Fluorobiphenyl	0.400	0.386	3.5	66	-0.01
16		Acenaphthylene	0.400	0.376	6.0	64	-0.01
17		Acenaphthene	0.400	0.381	4.8	64	-0.01
18		Fluorene	0.400	0.415	-3.7	71	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	83	0.00
20		4-Bromophenyl-phenylether	0.400	0.369	7.8	73	-0.01
21		Hexachlorobenzene	0.400	0.397	0.8	78	-0.01
22		Pentachlorophenol	0.400	0.021	94.8#	5	0.00
23		Phenanthrene	0.400	0.380	5.0	78	-0.01
24		Anthracene	0.400	0.355	11.3	74	-0.01
25	SURR	Fluoranthene-d10	0.400	0.407	-1.7	83	-0.01
26		Fluoranthene	0.400	0.412	-3.0	83	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	77	-0.01
28		Pyrene	0.400	0.442	-10.5	83	-0.01
29	S	Terphenyl-d14	0.400	0.417	-4.2	79	-0.02
30		Benzo(a)anthracene	0.400	0.363	9.3	68	0.00
31		Chrysene	0.400	0.399	0.3	74	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.377	5.8	71	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.355	11.3	63	0.00
34	I	Perylene-d12	0.400	0.400	0.0	72	0.00
35		Benzo(b)fluoranthene	0.400	0.358	10.5	64	-0.01
36		Benzo(k)fluoranthene	0.400	0.434	-8.5	75	-0.01
37	C	Benzo(a)pyrene	0.400	0.384	4.0	66	0.00
38		Dibenzo(a,h)anthracene	0.400	0.342	14.5	61	0.00
39		Benzo(g,h,i)perylene	0.400	0.415	-3.7	72	-0.01

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

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