Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060619\
Data File : BE099843.D
Acq On : 7 Jun 2019 4:03
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 07 05:45:00 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 05 14:35:58 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	103	0.00
2		1,4-Dioxane	0.400	0.279	30.3#	84	0.01
3		n-Nitrosodimethylamine	0.400	0.354	11.5	111	0.01
4	S	2-Fluorophenol	0.400	0.410	-2.5	123	0.01
5	S	Phenol-d6	0.400	0.480	-20.0	138	0.01
6		bis(2-Chloroethyl)ether	0.400	0.377	5.8	113	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	109	0.00
8	S	Nitrobenzene-d5	0.400	0.405	-1.3	126	0.00
9		Naphthalene	0.400	0.392	2.0	119	0.01
10		Hexachlorobutadiene	0.400	0.371	7.3	113	0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.383	4.3	117	0.01
12		2-Methylnaphthalene	0.400	0.408	-2.0	127	0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	111	0.01
14	S	2,4,6-Tribromophenol	0.400	0.543	-35.8#	177	0.00
15	S	2-Fluorobiphenyl	0.400	0.362	9.5	117	0.00
16		Acenaphthylene	0.400	0.400	0.0	126	0.00
17		Acenaphthene	0.400	0.420	-5.0	132	0.00
18		Fluorene	0.400	0.438	-9.5	133	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	90	0.00
20		4-Bromophenyl-phenylether	0.400	0.398	0.5	110	0.01
21		Hexachlorobenzene	0.400	0.350	12.5	92	0.01
22		Pentachlorophenol	0.400	0.806	-101.5#	305	-0.03
23		Phenanthrene	0.400	0.465	-16.3	119	0.00
24		Anthracene	0.400	0.468	-17.0	126	0.00
25	SURR	Fluoranthene-d10	0.400	0.305	23.8	79	0.00
26		Fluoranthene	0.400	0.392	2.0	100	0.01
27	I	Chrysene-d12	0.400	0.400	0.0	72	0.00
28		Pyrene	0.400	0.526	-31.5#	97	0.01
29	S	Terphenyl-d14	0.400	0.477	-19.2	88	0.00
30		Benzo(a)anthracene	0.400	0.508	-27.0#	96	0.01
31		Chrysene	0.400	0.402	-0.5	73	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.660	-65.0#	123	0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.275	31.3#	51	0.05
34	I	Perylene-d12	0.400	0.400	0.0	70	0.03
35		Benzo(b)fluoranthene	0.400	0.406	-1.5	79	0.02
36		Benzo(k)fluoranthene	0.400	0.356	11.0	69	0.02
37	C	Benzo(a)pyrene	0.400	0.401	-0.3	78	0.02
38		Dibenzo(a,h)anthracene	0.400	0.282	29.5#	56	0.04
39		Benzo(g,h,i)perylene	0.400	0.266	33.5#	53	0.06

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

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