Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE070219\
Data File : BE100293.D
Acq On : 2 Jul 2019 9:35
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 03 02:32:27 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	101	0.00
2		1,4-Dioxane	0.400	0.366	8.5	89	0.00
3		n-Nitrosodimethylamine	0.400	0.454	-13.5	139	-0.01
4	S	2-Fluorophenol	0.400	0.535	-33.8#	134	0.00
5	S	Phenol-d6	0.400	0.442	-10.5	112	0.00
6		bis(2-Chloroethyl)ether	0.400	0.424	-6.0	96	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	91	-0.01
8	S	Nitrobenzene-d5	0.400	0.442	-10.5	103	-0.01
9		Naphthalene	0.400	0.420	-5.0	94	0.00
10		Hexachlorobutadiene	0.400	0.402	-0.5	89	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	91	0.00
12		2-Methylnaphthalene	0.400	0.390	2.5	93	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	93	0.02
14	S	2,4,6-Tribromophenol	0.400	0.613	-53.2#	145	0.00
15	S	2-Fluorobiphenyl	0.400	0.419	-4.7	95	0.00
16		Acenaphthylene	0.400	0.446	-11.5	100	0.00
17		Acenaphthene	0.400	0.427	-6.7	94	0.00
18		Fluorene	0.400	0.434	-8.5	98	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	92	0.01
20		4-Bromophenyl-phenylether	0.400	0.432	-8.0	95	-0.01
21		Hexachlorobenzene	0.400	0.426	-6.5	94	0.01
22		Pentachlorophenol	0.400	0.124	69.0#	32	0.01
23		Phenanthrene	0.400	0.440	-10.0	101	0.00
24		Anthracene	0.400	0.470	-17.5	110	0.00
25	SURR	Fluoranthene-d10	0.400	0.433	-8.2	98	0.00
26		Fluoranthene	0.400	0.460	-15.0	103	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	89	0.00
28		Pyrene	0.400	0.491	-22.7	107	0.00
29	S	Terphenyl-d14	0.400	0.442	-10.5	97	0.00
30		Benzo(a)anthracene	0.400	0.434	-8.5	95	0.01
31		Chrysene	0.400	0.455	-13.7	99	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.462	-15.5	98	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.410	-2.5	84	0.02
34	I	Perylene-d12	0.400	0.400	0.0	85	0.01
35		Benzo(b)fluoranthene	0.400	0.441	-10.2	93	0.00
36		Benzo(k)fluoranthene	0.400	0.402	-0.5	82	0.01
37	C	Benzo(a)pyrene	0.400	0.434	-8.5	89	0.02
38		Dibenzo(a,h)anthracene	0.400	0.411	-2.7	87	0.02
39		Benzo(g,h,i)perylene	0.400	0.413	-3.2	85	0.02

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

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