Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE101918\
Data File : BE098001.D
Acq On : 19 Oct 2018 20:56
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 19 23:21:07 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 16 15:36:44 2018
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	-0.01
2		1,4-Dioxane	0.400	0.436	-9.0	120	0.00
3		n-Nitrosodimethylamine	0.400	0.382	4.5	109	0.00
4	S	2-Fluorophenol	0.400	0.430	-7.5	115	-0.01
5	S	Phenol-d6	0.400	0.477	-19.2	124	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.405	-1.3	110	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	118	-0.01
8	S	Nitrobenzene-d5	0.400	0.420	-5.0	127	-0.01
9		Naphthalene	0.400	0.434	-8.5	132	-0.01
10		Hexachlorobutadiene	0.400	0.449	-12.2	137	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.461	-15.3	140	-0.02
12		2-Methylnaphthalene	0.400	0.453	-13.2	139	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	129	-0.02
14	S	2,4,6-Tribromophenol	0.400	0.389	2.8	142	-0.01
15	S	2-Fluorobiphenyl	0.400	0.401	-0.3	137	-0.02
16		Acenaphthylene	0.400	0.420	-5.0	143	0.00
17		Acenaphthene	0.400	0.420	-5.0	142	0.00
18		Fluorene	0.400	0.408	-2.0	138	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	122	-0.01
20		4-Bromophenyl-phenylether	0.400	0.443	-10.7	139	-0.01
21		Hexachlorobenzene	0.400	0.463	-15.8	145	0.00
22		Pentachlorophenol	0.400	0.589	-47.2#	214	-0.01
23		Phenanthrene	0.400	0.411	-2.7	130	-0.01
24		Anthracene	0.400	0.412	-3.0	132	-0.01
25	SURR	Fluoranthene-d10	0.400	0.384	4.0	126	-0.01
26		Fluoranthene	0.400	0.384	4.0	126	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	111	-0.01
28		Pyrene	0.400	0.420	-5.0	124	-0.01
29	S	Terphenyl-d14	0.400	0.415	-3.7	127	-0.01
30		Benzo(a)anthracene	0.400	0.418	-4.5	123	0.00
31		Chrysene	0.400	0.434	-8.5	131	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.397	0.8	113	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.426	-6.5	120	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	109	-0.02
35		Benzo(b)fluoranthene	0.400	0.418	-4.5	122	-0.01
36		Benzo(k)fluoranthene	0.400	0.427	-6.7	128	-0.01
37	C	Benzo(a)pyrene	0.400	0.425	-6.2	123	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.414	-3.5	120	-0.02
39		Benzo(g,h,i)perylene	0.400	0.406	-1.5	121	-0.02

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

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