Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN101718\
Data File : BN003181.D
Acq On : 17 Oct 2018 15:50
Operator : MJ/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Oct 18 15:50:16 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN101718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Oct 18 15:26:23 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	52	0.02
2		1,4-Dioxane	0.400	0.392	2.0	57	0.00
3		n-Nitrosodimethylamine	0.400	0.211	47.3#	13	0.03
4	S	2-Fluorophenol	0.400	0.322	19.5	40	0.04
5	S	Phenol-d6	0.400	0.324	19.0	45	0.07
6		bis(2-Chloroethyl)ether	0.400	0.610	-52.5#	76	0.05
7	I	Naphthalene-d8	0.400	0.400	0.0	58	0.04
8	S	Nitrobenzene-d5	0.400	0.347	13.3	55	0.08
9		Naphthalene	0.400	0.373	6.8	56	0.04
10		Hexachlorobutadiene	0.400	0.383	4.3	57	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.369	7.8	56	0.05
12		2-Methylnaphthalene	0.400	0.361	9.8	55	0.04
13	I	Acenaphthene-d10	0.400	0.400	0.0	63	0.01
14	S	2,4,6-Tribromophenol	0.400	0.321	19.8	58	0.02
15	S	2-Fluorobiphenyl	0.400	0.338	15.5	56	0.03
16		Acenaphthylene	0.400	0.342	14.5	56	0.02
17		Acenaphthene	0.400	0.366	8.5	60	0.01
18		Fluorene	0.400	0.359	10.3	59	0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	65	0.02
20		4-Bromophenyl-phenylether	0.400	0.324	19.0	56	0.02
21		Hexachlorobenzene	0.400	0.351	12.3	59	0.01
22		Pentachlorophenol	0.400	0.353	11.8	63	0.04
23		Phenanthrene	0.400	0.352	12.0	60	0.01
24		Anthracene	0.400	0.340	15.0	58	0.03
25	SURR	Fluoranthene-d10	0.400	0.372	7.0	63	0.01
26		Fluoranthene	0.400	0.371	7.3	63	0.02
27	I	Chrysene-d12	0.400	0.400	0.0	72	0.01
28		Pyrene	0.400	0.329	17.8	64	0.01
29	S	Terphenyl-d14	0.400	0.322	19.5	64	0.01
30		Benzo(a)anthracene	0.400	0.320	20.0	66	0.01
31		Chrysene	0.400	0.368	8.0	68	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.320	20.0	61	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.391	2.3	77	0.02
34	I	Perylene-d12	0.400	0.400	0.0	78	0.01
35		Benzo(b)fluoranthene	0.400	0.338	15.5	72	0.00
36		Benzo(k)fluoranthene	0.400	0.364	9.0	74	0.00
37	C	Benzo(a)pyrene	0.400	0.355	11.3	74	0.01
38		Dibenzo(a,h)anthracene	0.400	0.369	7.8	77	0.02
39		Benzo(g,h,i)perylene	0.400	0.373	6.8	78	0.01

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

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