Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\Data\Be051816\
Data File : BE091791.D
Acq On : 18 May 2016 13:49
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 18 18:03:23 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 26 16:09:46 2016
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	5.000	5.000	0.0	89	-0.02
2		1,4-Dioxane	0.400	0.302	24.5	67	-0.02
3		n-Nitrosodimethylamine	0.400	0.329	17.8	74	0.00
4	S	2-Fluorophenol	0.400	0.356	11.0	82	-0.02
5	S	Phenol-d6	0.400	0.352	12.0	83	0.00
6		bis(2-Chloroethyl)ether	0.400	0.379	5.3	83	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	91	-0.02
8	S	Nitrobenzene-d5	0.400	0.350	12.5	85	-0.01
9		Nitrobenzene	0.400	0.336	16.0	83	-0.01
10		Naphthalene	0.400	0.382	4.5	86	0.00
11		Hexachlorobutadiene	0.400	0.383	4.3	89	-0.02
12		2-Methylnaphthalene	0.400	0.379	5.3	86	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	94	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.489	-22.2	102	-0.01
15	S	2-Fluorobiphenyl	0.400	0.373	6.8	88	-0.01
16		Acenaphthylene	0.400	0.394	1.5	105	0.00
17		Acenaphthene	0.400	0.358	10.5	84	-0.01
18		Fluorene	0.400	0.389	2.8	91	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	98	-0.01
20		4-Bromophenyl-phenylether	0.400	0.383	4.3	93	0.00
21		Hexachlorobenzene	0.400	0.389	2.8	95	0.00
22		Pentachlorophenol	0.400	0.388	3.0	97	0.00
23		Phenanthrene	0.400	0.397	0.8	95	-0.01
24		Anthracene	0.400	0.382	4.5	95	-0.01
25		Fluoranthene	0.400	0.429	-7.2	104	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	121	-0.02
27		Benzidine	0.400	0.354	11.5	102	0.00
28		Pyrene	0.400	0.316	21.0	99	0.00
29	S	Terphenyl-d14	0.400	0.300	25.0#	93	-0.02
30		Benzo(a)anthracene	0.400	0.332	17.0	103	-0.02
31		3,3'-Dichlorobenzidine	0.400	0.374	6.5	131	-0.02
32		Chrysene	0.400	0.300	25.0#	86	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.451	-12.7	177	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.397	0.8	118	-0.03
35	I	Perylene-d12	5.000	5.000	0.0	123	-0.02
36		Benzo(b)fluoranthene	0.400	0.349	12.8	111	-0.02
37		Benzo(k)fluoranthene	0.400	0.354	11.5	114	-0.02
38	C	Benzo(a)pyrene	0.400	0.358	10.5	117	-0.02
39		Dibenzo(a,h)anthracene	0.400	0.367	8.3	116	-0.05
40		Benzo(g,h,i)perylene	0.400	0.372	7.0	119	-0.05

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

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