Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051917\
Data File : BE093055.D
Acq On : 19 May 2017 11:28
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 19 17:51:44 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 15 13:35:57 2017
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	72	0.00
2		1,4-Dioxane	0.400	0.414	-3.5	85	0.00
3		n-Nitrosodimethylamine	0.400	0.424	-6.0	87	0.00
4	S	2-Fluorophenol	0.400	0.420	-5.0	85	0.00
5	S	Phenol-d6	0.400	0.432	-8.0	89	0.00
6		bis(2-Chloroethyl)ether	0.400	0.414	-3.5	85	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	76	0.00
8	S	Nitrobenzene-d5	0.400	0.423	-5.7	93	0.00
9		Naphthalene	0.400	0.390	2.5	83	0.00
10		Hexachlorobutadiene	0.400	0.391	2.3	84	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	86	0.00
12		2-Methylnaphthalene	0.400	0.400	0.0	86	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	80	0.00
14	S	2,4,6-Tribromophenol	0.400	0.448	-12.0	104	0.00
15	S	2-Fluorobiphenyl	0.400	0.392	2.0	87	0.00
16		Acenaphthylene	0.400	0.396	1.0	93	0.00
17		Acenaphthene	0.400	0.398	0.5	90	0.00
18		Fluorene	0.400	0.413	-3.2	93	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	79	0.00
20		4-Bromophenyl-phenylether	0.400	0.374	6.5	93	0.00
21		Hexachlorobenzene	0.400	0.354	11.5	99	0.00
22		Pentachlorophenol	0.400	0.359	10.3	102	0.00
23		Phenanthrene	0.400	0.384	4.0	95	0.03
24		Anthracene	0.400	0.388	3.0	107	0.00
25	SURR	Fluoranthene-d10	0.400	0.416	-4.0	102	0.00
26		Fluoranthene	0.400	0.414	-3.5	104	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	96	-0.02
28		Pyrene	0.400	0.402	-0.5	105	0.00
29	S	Terphenyl-d14	0.400	0.382	4.5	100	0.00
30		Benzo(a)anthracene	0.400	0.406	-1.5	109	0.00
31		Chrysene	0.400	0.389	2.8	105	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.474	-18.5	142	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.384	4.0	104	0.00
34	I	Perylene-d12	0.400	0.400	0.0	96	-0.01
35		Benzo(b)fluoranthene	0.400	0.399	0.3	112	0.00
36		Benzo(k)fluoranthene	0.400	0.375	6.3	102	0.00
37	C	Benzo(a)pyrene	0.400	0.387	3.3	109	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.379	5.3	105	0.00
39		Benzo(g,h,i)perylene	0.400	0.389	2.8	107	0.00

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

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