Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051916\
Data File : BE091809.D
Acq On : 19 May 2016 14:26
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE051916

Quant Time: May 19 15:43:39 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 19 15:36:49 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	99	-0.02
2		1,4-Dioxane	0.400	0.445	-11.2	93	0.00
3		n-Nitrosodimethylamine	0.400	0.355	11.3	96	0.00
4	S	2-Fluorophenol	0.400	0.360	10.0	95	0.00
5	S	Phenol-d6	0.400	0.340	15.0	94	0.00
6		bis(2-Chloroethyl)ether	0.400	0.354	11.5	96	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	97	0.00
8	S	Nitrobenzene-d5	0.400	0.346	13.5	94	0.00
9		Nitrobenzene	0.400	0.333	16.8	93	0.00
10		Naphthalene	0.400	0.380	5.0	96	0.00
11		Hexachlorobutadiene	0.400	0.385	3.8	97	-0.02
12		2-Methylnaphthalene	0.400	0.371	7.3	95	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	95	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.330	17.5	91	-0.01
15	S	2-Fluorobiphenyl	0.400	0.378	5.5	95	0.00
16		Acenaphthylene	0.400	0.371	7.3	93	0.00
17		Acenaphthene	0.400	0.367	8.3	93	-0.01
18		Fluorene	0.400	0.375	6.3	96	-0.01
19	I	Phenanthrene-d10	5.000	5.000	0.0	99	-0.01
20		4-Bromophenyl-phenylether	0.400	0.381	4.8	96	0.00
21		Hexachlorobenzene	0.400	0.387	3.3	97	0.00
22		Pentachlorophenol	0.400	0.317	20.8	87	0.00
23		Phenanthrene	0.400	0.386	3.5	97	-0.01
24		Anthracene	0.400	0.372	7.0	95	-0.01
25		Fluoranthene	0.400	0.379	5.3	94	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	101	-0.02
27		Benzidine	0.400	0.327	18.3	95	0.00
28		Pyrene	0.400	0.359	10.3	95	0.00
29	S	Terphenyl-d14	0.400	0.374	6.5	101	-0.02
30		Benzo(a)anthracene	0.400	0.375	6.3	101	-0.02
31		3,3'-Dichlorobenzidine	0.400	0.316	21.0	91	-0.02
32		Chrysene	0.400	0.437	-9.2	102	0.00
33		Bis(2-ethylhexyl)phthalate	0.400	0.453	-13.2	88	-0.02
34		Indeno(1,2,3-cd)pyrene	0.400	0.356	11.0	95	-0.03
35	I	Perylene-d12	5.000	5.000	0.0	99	-0.02
36		Benzo(b)fluoranthene	0.400	0.364	9.0	95	-0.02
37		Benzo(k)fluoranthene	0.400	0.369	7.8	96	-0.02
38	C	Benzo(a)pyrene	0.400	0.362	9.5	95	-0.02
39		Dibenzo(a,h)anthracene	0.400	0.361	9.8	95	-0.05
40		Benzo(g,h,i)perylene	0.400	0.369	7.8	96	-0.05

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

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