Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE052516\
Data File : BE091864.D
Acq On : 25 May 2016 18:30
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE052516

Quant Time: May 25 19:47:24 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE052516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 25 19:42:30 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	97	0.00
2		1,4-Dioxane	0.400	0.418	-4.5	91	0.00
3		n-Nitrosodimethylamine	0.400	0.319	20.3	88	0.00
4	S	2-Fluorophenol	0.400	0.346	13.5	90	0.00
5	S	Phenol-d6	0.400	0.326	18.5	89	0.00
6		bis(2-Chloroethyl)ether	0.400	0.342	14.5	90	-0.02
7		n-Nitroso-di-n-propylamine	0.400	0.327	18.3	92	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	96	0.00
9	S	Nitrobenzene-d5	0.400	0.345	13.8	91	0.00
10		Nitrobenzene	0.400	0.327	18.3	90	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.440	-10.0	134	0.00
12		Naphthalene	0.400	0.363	9.3	93	0.00
13		Hexachlorobutadiene	0.400	0.369	7.8	92	0.00
14		2-Methylnaphthalene	0.400	0.359	10.3	93	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
16	S	2,4,6-Tribromophenol	0.400	0.321	19.8	90	0.00
17	S	2-Fluorobiphenyl	0.400	0.357	10.8	92	0.00
18		Acenaphthylene	0.400	0.347	13.3	91	0.00
19		2,6-Dinitrotoluene	0.400	0.404	-1.0	121	0.00
20		Acenaphthene	0.400	0.354	11.5	92	0.02
21		2,4-Dinitrotoluene	0.400	0.423	-5.7	89	0.02
22		Fluorene	0.400	0.348	13.0	92	0.00
23		4-Chlorophenyl-phenylether	0.400	0.328	18.0	93	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
25		4-Bromophenyl-phenylether	0.400	0.354	11.5	93	0.00
26		Hexachlorobenzene	0.400	0.359	10.3	91	0.00
27		Pentachlorophenol	0.400	0.329	17.8	104	0.00
28		Phenanthrene	0.400	0.335	16.3	92	0.00
29		Anthracene	0.400	0.343	14.2	92	0.00
30		Fluoranthene	0.400	0.325	18.8	91	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	90	0.00
32		Benzidine	0.400	0.378	5.5	87	0.02
33		Pyrene	0.400	0.381	4.8	93	0.00
34	S	Terphenyl-d14	0.400	0.378	5.5	90	0.00
35		Benzo(a)anthracene	0.400	0.374	6.5	89	0.00
36		3,3'-Dichlorobenzidine	0.400	0.348	13.0	86	0.00
37		Chrysene	0.400	0.332	17.0	93	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.359	10.3	94	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.376	6.0	90	0.00
40	I	Perylene-d12	5.000	5.000	0.0	97	0.00
41		Benzo(b)fluoranthene	0.400	0.364	9.0	101	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.334	16.5	79	0.00
43 C	Benzo(a)pyrene	0.400	0.343	14.2	89	0.00
44	Dibenzo(a,h)anthracene	0.400	0.345	13.8	89	0.00
45	Benzo(g,h,i)perylene	0.400	0.352	12.0	91	0.00

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