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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
2		1,4-Dioxane	0.737	0.647	12.2	91	0.00
3		n-Nitrosodimethylamine	0.918	0.732	20.3	88	0.00
4	S	2-Fluorophenol	1.429	1.235	13.6	90	0.00
5	S	Phenol-d6	1.874	1.526	18.6	89	0.00
6		bis(2-Chloroethyl)ether	1.569	1.340	14.6	90	-0.02
7		n-Nitroso-di-n-propylamine	1.452	1.189	18.1	92	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
9	S	Nitrobenzene-d5	0.406	0.351	13.5	91	0.00
10		Nitrobenzene	0.394	0.322	18.3	90	0.00
11		bis(2-Chloroethoxy)methane	0.474	0.521	-9.9	134	0.00
12		Naphthalene	1.080	0.980	9.3	93	0.00
13		Hexachlorobutadiene	0.166	0.153	7.8	92	0.00
14		2-Methylnaphthalene	0.721	0.647	10.3	93	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
16	S	2,4,6-Tribromophenol	0.223	0.179	19.7	90	0.00
17	S	2-Fluorobiphenyl	1.487	1.325	10.9	92	0.00
18		Acenaphthylene	2.458	2.046	16.8	91	0.00
19		2,6-Dinitrotoluene	0.341	0.345	-1.2	121	0.00
20		Acenaphthene	1.337	1.162	13.1	92	0.02
21		2,4-Dinitrotoluene	0.393	0.263	33.1#	89	0.02
22		Fluorene	1.755	1.525	13.1	92	0.00
23		4-Chlorophenyl-phenylether	0.883	0.780	11.7	93	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
25		4-Bromophenyl-phenylether	0.179	0.158	11.7	93	0.00
26		Hexachlorobenzene	0.170	0.153	10.0	91	0.00
27		Pentachlorophenol	0.066	0.054	18.2	104	0.00
28		Phenanthrene	1.090	0.954	12.5	92	0.00
29		Anthracene	0.999	0.858	14.1	92	0.00
30		Fluoranthene	1.273	1.090	14.4	91	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	90	0.00
32		Benzidine	0.400	0.313	21.8	87	0.02
33		Pyrene	1.061	1.010	4.8	93	0.00
34	S	Terphenyl-d14	0.755	0.714	5.4	90	0.00
35		Benzo(a)anthracene	1.333	1.246	6.5	89	0.00
36		3,3'-Dichlorobenzidine	0.732	0.638	12.8	86	0.00
37		Chrysene	1.222	1.237	-1.2	93	0.00
38		Bis(2-ethylhexyl)phthalate	0.484	0.434	10.3	94	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.176	5.9	90	0.00
40	I	Perylene-d12	1.000	1.000	0.0	97	0.00
41		Benzo(b)fluoranthene	1.349	1.228	9.0	101	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.225	1.022	16.6	79	0.00
43 C	Benzo(a)pyrene	1.223	1.048	14.3	89	0.00
44	Dibenzo(a,h)anthracene	1.172	1.011	13.7	89	0.00
45	Benzo(g,h,i)perylene	1.189	1.046	12.0	91	0.00

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