Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072716\
Data File : BE092058.D
Acq On : 27 Jul 2016 15:52
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE072716

Quant Time: Jul 27 17:39:13 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 27 17:13:26 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	92	-0.02
2		1,4-Dioxane	0.687	0.704	-2.5	95	0.00
3		n-Nitrosodimethylamine	0.819	0.749	8.5	93	0.00
4	S	2-Fluorophenol	1.114	0.992	11.0	94	-0.02
5	S	Phenol-d6	1.487	1.206	18.9	85	0.00
6		bis(2-Chloroethyl)ether	1.437	1.292	10.1	92	0.00
7		n-Nitroso-di-n-propylamine	1.187	1.058	10.9	93	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
9	S	Nitrobenzene-d5	0.373	0.305	18.2	81	0.00
10		Nitrobenzene	0.362	0.317	12.4	91	0.00
11		bis(2-Chloroethoxy)methane	0.417	0.362	13.2	92	0.00
12		Naphthalene	1.095	1.075	1.8	91	0.00
13		Hexachlorobutadiene	0.164	0.164	0.0	94	0.00
14		2-Methylnaphthalene	0.707	0.667	5.7	92	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
16	S	2,4,6-Tribromophenol	0.175	0.147	16.0	89	0.00
17	S	2-Fluorobiphenyl	2.003	2.048	-2.2	91	0.00
18		Acenaphthylene	12.962	12.552	3.2	92	0.00
19		2,6-Dinitrotoluene	1.331	1.158	13.0	87	0.00
20		Acenaphthene	1.366	1.363	0.2	89	0.00
21		2,4-Dinitrotoluene	1.684	0.969	42.5#	81	0.00
22		Fluorene	2.145	2.110	1.6	89	0.00
23		4-Chlorophenyl-phenylether	1.055	1.104	-4.6	92	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
25		4-Bromophenyl-phenylether	0.206	0.198	3.9	88	0.00
26		Hexachlorobenzene	0.213	0.210	1.4	89	0.00
27		Pentachlorophenol	0.060	0.030	50.0#	82	0.00
28		Phenanthrene	1.271	1.257	1.1	91	0.00
29		Anthracene	1.156	1.069	7.5	88	0.00
30		Fluoranthene	5.652	5.158	8.7	92	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
32		Benzidine	0.246	0.190	22.8	85	0.00
33		Pyrene	6.442	6.683	-3.7	90	0.00
34	S	Terphenyl-d14	0.931	0.934	-0.3	89	0.00
35		Benzo(a)anthracene	2.391	2.412	-0.9	86	0.00
36		3,3'-Dichlorobenzidine	0.461	0.349	24.3	90	0.00
37		Chrysene	2.026	2.079	-2.6	79	0.00
38		Bis(2-ethylhexyl)phthalate	4.801	5.020	-4.6	98	0.00
39		Indeno(1,2,3-cd)pyrene	5.859	6.595	-12.6	96	0.00
40	I	Perylene-d12	1.000	1.000	0.0	93	0.00
41		Benzo(b)fluoranthene	1.515	1.521	-0.4	97	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.567	1.557	0.6	91	0.00
43 C	Benzo(a)pyrene	1.273	1.198	5.9	90	0.00
44	Dibenzo(a,h)anthracene	1.111	1.062	4.4	96	0.00
45	Benzo(g,h,i)perylene	4.638	4.521	2.5	97	0.00

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