Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be072216\
Data File : BE092007.D
Acq On : 22 Jul 2016 15:29
Operator : UM/SJ
Sample : SSTDICV0.4
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE072216

Quant Time: Jul 22 16:31:32 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 22 15:39:27 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	94	0.00
2		1,4-Dioxane	0.676	0.731	-8.1	96	-0.02
3		n-Nitrosodimethylamine	0.753	0.786	-4.4	108	-0.02
4	S	2-Fluorophenol	0.969	1.034	-6.7	110	0.00
5	S	Phenol-d6	1.312	1.381	-5.3	108	-0.02
6		bis(2-Chloroethyl)ether	1.436	1.475	-2.7	105	-0.02
7		n-Nitroso-di-n-propylamine	1.226	1.181	3.7	100	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
9	S	Nitrobenzene-d5	0.301	0.307	-2.0	101	-0.02
10		Nitrobenzene	0.341	0.327	4.1	100	-0.02
11		bis(2-Chloroethoxy)methane	0.377	0.370	1.9	101	-0.03
12		Naphthalene	1.136	1.130	0.5	92	0.00
13		Hexachlorobutadiene	0.163	0.164	-0.6	92	0.00
14		2-Methylnaphthalene	0.712	0.710	0.3	95	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	81	0.00
16	S	2,4,6-Tribromophenol	0.122	0.123	-0.8	107	0.00
17	S	2-Fluorobiphenyl	1.315	1.477	-12.3	95	0.00
18		Acenaphthylene	7.691	8.751	-13.8	113	0.00
19		2,6-Dinitrotoluene	0.926	0.968	-4.5	125	0.00
20		Acenaphthene	1.304	1.265	3.0	79	0.00
21		2,4-Dinitrotoluene	0.912	0.790	13.4	105	0.00
22		Fluorene	1.556	1.756	-12.9	99	0.00
23		4-Chlorophenyl-phenylether	0.772	0.901	-16.7	98	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
25		4-Bromophenyl-phenylether	0.208	0.210	-1.0	100	0.00
26		Hexachlorobenzene	0.215	0.218	-1.4	97	0.00
27		Pentachlorophenol	0.053	0.036	32.1#	115	0.00
28		Phenanthrene	1.292	1.322	-2.3	99	-0.01
29		Anthracene	1.171	1.183	-1.0	104	-0.01
30		Fluoranthene	5.165	4.968	3.8	100	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	87	0.00
32		Benzidine	0.156	0.178	-14.1	125	0.00
33		Pyrene	3.628	3.700	-2.0	95	0.00
34	S	Terphenyl-d14	0.502	0.566	-12.7	102	0.00
35		Benzo(a)anthracene	1.127	1.239	-9.9	107	0.00
36		3,3'-Dichlorobenzidine	0.169	0.165	2.4	112	0.00
37		Chrysene	1.372	1.572	-14.6	100	0.00
38		Bis(2-ethylhexyl)phthalate	0.811	0.915	-12.8	119	0.00
39		Indeno(1,2,3-cd)pyrene	3.769	3.723	1.2	87	0.00
40	I	Perylene-d12	1.000	1.000	0.0	93	0.00
41		Benzo(b)fluoranthene	1.312	1.260	4.0	94	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.337	1.305	2.4	90	0.00
43 C	Benzo(a)pyrene	1.221	1.139	6.7	90	0.00
44	Dibenzo(a,h)anthracene	1.089	1.008	7.4	88	0.00
45	Benzo(g,h,i)perylene	4.463	4.292	3.8	92	0.00

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