Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
Data File : BE090369.D
Acq On : 5 Aug 2015 21:07
Operator : TP/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE080515

Quant Time: Aug 06 13:01:22 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 2015
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	102	0.00
2		1,4-Dioxane	0.687	0.785	-14.3	109	0.00
3		n-Nitrosodimethylamine	0.773	0.715	7.5	95	0.00
4	S	2-Fluorophenol	0.750	0.704	6.1	103	0.00
5	S	Phenol-d6	1.241	1.197	3.5	106	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
7	S	Nitrobenzene-d5	0.331	0.329	0.6	108	0.00
8		Nitrobenzene	0.349	0.336	3.7	110	0.00
9		Naphthalene	1.104	1.108	-0.4	104	0.00
10		Hexachlorobutadiene	0.191	0.190	0.5	101	0.00
11		2-Methylnaphthalene	0.640	0.665	-3.9	111	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
13	S	2,4,6-Tribromophenol	0.095	0.083	12.6	118	0.00
14	S	2-Fluorobiphenyl	1.375	1.447	-5.2	109	0.00
15		Acenaphthylene	8.098	8.418	-4.0	111	0.00
16		Acenaphthene	1.259	1.261	-0.2	105	0.00
17		Fluorene	1.635	1.707	-4.4	109	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
19		4-Bromophenyl-phenylether	0.194	0.204	-5.2	110	0.00
20		Hexachlorobenzene	1.005	1.042	-3.7	106	0.00
21		Pentachlorophenol	0.044	0.032	27.3#	120	0.00
22		Phenanthrene	1.251	1.233	1.4	102	0.00
23		Anthracene	1.028	1.071	-4.2	110	0.00
24		Fluoranthene	1.231	1.280	-4.0	111	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
26		Pyrene	1.084	1.083	0.1	110	0.00
27	S	Terphenyl-d14	0.737	0.743	-0.8	107	0.00
28		Benzo(a)anthracene	1.108	1.050	5.2	111	0.00
29		Chrysene	1.339	1.386	-3.5	107	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.224	28.0#	105	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	0.959	-1.8	130	0.00
32	I	Perylene-d12	1.000	1.000	0.0	107	0.00
33		Benzo(b)fluoranthene	0.918	0.966	-5.2	120	0.00
34		Benzo(k)fluoranthene	1.392	1.415	-1.7	106	0.00
35	C	Benzo(a)pyrene	0.850	0.890	-4.7	119	0.00
36		Dibenzo(a,h)anthracene	0.766	0.820	-7.0	133	0.00
37		Benzo(g,h,i)perylene	0.923	0.988	-7.0	119	0.00

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