Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090415\
Data File : BE090745.D
Acq On : 5 Sep 2015 00:29
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 07 01:03:25 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	78	0.00
2		1,4-Dioxane	0.814	0.768	5.7	76	0.00
3		n-Nitrosodimethylamine	1.033	0.888	14.0	71	0.00
4	S	2-Fluorophenol	1.039	1.059	-1.9	85	0.00
5	S	Phenol-d6	1.392	1.486	-6.8	88	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
7	S	Nitrobenzene-d5	0.330	0.324	1.8	83	0.00
8		Nitrobenzene	0.365	0.346	5.2	84	0.00
9		Naphthalene	1.091	1.096	-0.5	86	0.00
10		Hexachlorobutadiene	0.171	0.167	2.3	80	0.00
11		2-Methylnaphthalene	0.590	0.665	-12.7	99	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
13	S	2,4,6-Tribromophenol	0.143	0.148	-3.5	120	0.00
14	S	2-Fluorobiphenyl	1.385	1.353	2.3	97	0.00
15		Acenaphthylene	7.992	8.412	-5.3	107	0.00
16		Acenaphthene	1.335	1.367	-2.4	99	0.00
17		Fluorene	1.666	1.730	-3.8	103	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
19		4-Bromophenyl-phenylether	0.170	0.177	-4.1	103	0.00
20		Hexachlorobenzene	0.928	0.911	1.8	95	0.00
21		Pentachlorophenol	0.054	0.053	1.9	106	0.00
22		Phenanthrene	1.237	1.272	-2.8	98	-0.02
23		Anthracene	1.035	1.082	-4.5	103	0.00
24		Fluoranthene	1.304	1.378	-5.7	108	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	96	0.00
26		Pyrene	1.035	1.158	-11.9	113	0.00
27	S	Terphenyl-d14	0.553	0.596	-7.8	108	0.00
28		Benzo(a)anthracene	1.156	1.181	-2.2	104	0.00
29		Chrysene	1.279	1.335	-4.4	101	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.335	2.9	127	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.337	-13.2	118	0.00
32	I	Perylene-d12	1.000	1.000	0.0	102	0.00
33		Benzo(b)fluoranthene	1.084	1.149	-6.0	111	0.00
34		Benzo(k)fluoranthene	1.263	1.310	-3.7	108	0.00
35	C	Benzo(a)pyrene	0.970	1.057	-9.0	119	0.00
36		Dibenzo(a,h)anthracene	1.020	1.110	-8.8	116	-0.01
37		Benzo(g,h,i)perylene	1.116	1.169	-4.7	113	-0.02

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