Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070615\
Data File : BE090060.D
Acq On : 6 Jul 2015 19:18
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 07 01:17:15 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	118	0.00
2		1,4-Dioxane	0.686	0.558	18.7	100	-0.01
3		n-Nitrosodimethylamine	0.930	0.758	18.5	100	-0.01
4	S	2-Fluorophenol	1.149	1.057	8.0	109	0.00
5	S	Phenol-d6	1.608	1.528	5.0	114	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	121	-0.01
7	S	Nitrobenzene-d5	0.397	0.341	14.1	106	0.00
8		Nitrobenzene	0.421	0.360	14.5	106	0.00
9		Naphthalene	1.130	1.041	7.9	113	-0.01
10		Hexachlorobutadiene	0.178	0.168	5.6	116	0.00
11		2-Methylnaphthalene	0.755	0.666	11.8	109	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	113	-0.01
13	S	2,4,6-Tribromophenol	0.147	0.116	21.1	91	0.00
14	S	2-Fluorobiphenyl	1.503	1.365	9.2	105	-0.01
15		Acenaphthylene	9.019	8.277	8.2	107	-0.01
16		Acenaphthene	1.300	1.212	6.8	106	-0.01
17		Fluorene	1.768	1.571	11.1	103	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.02
19		4-Bromophenyl-phenylether	0.213	0.185	13.1	92	-0.02
20		Hexachlorobenzene	0.885	0.792	10.5	97	-0.02
21		Pentachlorophenol	0.086	0.055	36.0#	70	0.00
22		Phenanthrene	1.281	1.118	12.7	93	-0.02
23		Anthracene	1.172	1.054	10.1	95	0.00
24		Fluoranthene	1.431	1.259	12.0	93	-0.01
25	I	Chrysene-d12	1.000	1.000	0.0	96	-0.01
26		Pyrene	1.192	1.134	4.9	93	-0.01
27	S	Terphenyl-d14	0.833	0.768	7.8	89	-0.01
28		Benzo(a)anthracene	1.249	1.148	8.1	91	-0.01
29		Chrysene	1.297	1.170	9.8	88	-0.01
30		Bis(2-ethylhexyl)phthalate	0.398	0.344	13.6	91	-0.01
31		Indeno(1,2,3-cd)pyrene	1.243	1.138	8.4	90	-0.04
32	I	Perylene-d12	1.000	1.000	0.0	91	-0.02
33		Benzo(b)fluoranthene	1.103	1.125	-2.0	94	-0.02
34		Benzo(k)fluoranthene	1.232	1.208	1.9	90	-0.02
35	C	Benzo(a)pyrene	1.019	1.030	-1.1	92	-0.03
36		Dibenzo(a,h)anthracene	1.046	0.994	5.0	88	-0.04
37		Benzo(g,h,i)perylene	1.060	1.005	5.2	87	-0.04

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