Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070915\
Data File : BE090076.D
Acq On : 9 Jul 2015 16:56
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 10 02:02:40 2015
Quant Method : Z:\HPCHEM1\BNA_E\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	114	-0.02
2		1,4-Dioxane	0.686	0.588	14.3	102	-0.01
3		n-Nitrosodimethylamine	0.930	0.746	19.8	95	-0.01
4	S	2-Fluorophenol	1.149	1.118	2.7	112	-0.01
5	S	Phenol-d6	1.608	1.732	-7.7	125	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	124	-0.02
7	S	Nitrobenzene-d5	0.397	0.350	11.8	111	-0.02
8		Nitrobenzene	0.421	0.365	13.3	111	-0.01
9		Naphthalene	1.130	1.040	8.0	116	-0.01
10		Hexachlorobutadiene	0.178	0.166	6.7	118	-0.01
11		2-Methylnaphthalene	0.755	0.694	8.1	116	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	120	-0.01
13	S	2,4,6-Tribromophenol	0.147	0.128	12.9	108	0.00
14	S	2-Fluorobiphenyl	1.503	1.380	8.2	113	-0.01
15		Acenaphthylene	9.019	8.254	8.5	113	0.00
16		Acenaphthene	1.300	1.210	6.9	113	-0.01
17		Fluorene	1.768	1.576	10.9	110	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	107	-0.02
19		4-Bromophenyl-phenylether	0.213	0.194	8.9	100	-0.02
20		Hexachlorobenzene	0.885	0.808	8.7	102	-0.02
21		Pentachlorophenol	0.086	0.144	-67.4#	191#	0.00
22		Phenanthrene	1.281	1.129	11.9	97	-0.02
23		Anthracene	1.172	1.044	10.9	97	-0.02
24		Fluoranthene	1.431	1.225	14.4	94	-0.01
25	I	Chrysene-d12	1.000	1.000	0.0	93	-0.01
26		Pyrene	1.192	1.188	0.3	94	-0.01
27	S	Terphenyl-d14	0.833	0.791	5.0	88	-0.01
28		Benzo(a)anthracene	1.249	1.161	7.0	88	-0.01
29		Chrysene	1.297	1.170	9.8	85	-0.01
30		Bis(2-ethylhexyl)phthalate	0.398	0.476	-19.6	121	-0.01
31		Indeno(1,2,3-cd)pyrene	1.243	1.196	3.8	91	-0.03
32	I	Perylene-d12	1.000	1.000	0.0	88	-0.02
33		Benzo(b)fluoranthene	1.103	1.153	-4.5	93	-0.02
34		Benzo(k)fluoranthene	1.232	1.234	-0.2	89	-0.02
35	C	Benzo(a)pyrene	1.019	1.055	-3.5	92	-0.02
36		Dibenzo(a,h)anthracene	1.046	1.030	1.5	89	-0.03
37		Benzo(g,h,i)perylene	1.060	1.051	0.8	89	-0.04

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