Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
Data File : BE090062.D
Acq On : 8 Jul 2015 16:51
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 09 01:19:06 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	138	-0.01
2		1,4-Dioxane	0.686	0.580	15.5	121	-0.01
3		n-Nitrosodimethylamine	0.930	0.766	17.6	118	-0.02
4	S	2-Fluorophenol	1.149	0.984	14.4	118	-0.01
5	S	Phenol-d6	1.608	1.448	10.0	126	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	140	-0.01
7	S	Nitrobenzene-d5	0.397	0.344	13.4	123	-0.01
8		Nitrobenzene	0.421	0.361	14.3	124	-0.01
9		Naphthalene	1.130	1.038	8.1	131	-0.01
10		Hexachlorobutadiene	0.178	0.166	6.7	133	-0.01
11		2-Methylnaphthalene	0.755	0.667	11.7	126	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	132	-0.01
13	S	2,4,6-Tribromophenol	0.147	0.098	33.3#	90	0.00
14	S	2-Fluorobiphenyl	1.503	1.356	9.8	122	-0.01
15		Acenaphthylene	9.019	8.220	8.9	124	-0.01
16		Acenaphthene	1.300	1.207	7.2	124	-0.01
17		Fluorene	1.768	1.578	10.7	121	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	122	-0.02
19		4-Bromophenyl-phenylether	0.213	0.187	12.2	110	-0.02
20		Hexachlorobenzene	0.885	0.785	11.3	113	-0.02
21		Pentachlorophenol	0.086	0.044	48.8#	67	0.00
22		Phenanthrene	1.281	1.120	12.6	110	-0.02
23		Anthracene	1.172	1.045	10.8	111	-0.02
24		Fluoranthene	1.431	1.249	12.7	109	-0.02
25	I	Chrysene-d12	1.000	1.000	0.0	113	-0.01
26		Pyrene	1.192	1.143	4.1	111	-0.02
27	S	Terphenyl-d14	0.833	0.762	8.5	104	-0.02
28		Benzo(a)anthracene	1.249	1.116	10.6	104	-0.01
29		Chrysene	1.297	1.169	9.9	104	-0.01
30		Bis(2-ethylhexyl)phthalate	0.398	0.242	39.2#	75	-0.01
31		Indeno(1,2,3-cd)pyrene	1.243	1.092	12.1	101	-0.04
32	I	Perylene-d12	1.000	1.000	0.0	104	-0.02
33		Benzo(b)fluoranthene	1.103	1.098	0.5	105	-0.02
34		Benzo(k)fluoranthene	1.232	1.245	-1.1	106	-0.02
35	C	Benzo(a)pyrene	1.019	1.000	1.9	103	-0.02
36		Dibenzo(a,h)anthracene	1.046	0.972	7.1	99	-0.03
37		Benzo(g,h,i)perylene	1.060	0.982	7.4	98	-0.04

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