Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070115\
Data File : BE090049.D
Acq On : 1 Jul 2015 20:12
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 02 05:16:31 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	102	0.00
2		1,4-Dioxane	0.686	0.589	14.1	91	0.00
3		n-Nitrosodimethylamine	0.930	0.768	17.4	88	-0.01
4	S	2-Fluorophenol	1.149	1.118	2.7	99	0.00
5	S	Phenol-d6	1.608	1.589	1.2	102	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	106	-0.02
7	S	Nitrobenzene-d5	0.397	0.346	12.8	94	0.00
8		Nitrobenzene	0.421	0.365	13.3	95	0.00
9		Naphthalene	1.130	1.040	8.0	99	0.00
10		Hexachlorobutadiene	0.178	0.166	6.7	100	0.00
11		2-Methylnaphthalene	0.755	0.677	10.3	96	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
13	S	2,4,6-Tribromophenol	0.147	0.129	12.2	89	0.00
14	S	2-Fluorobiphenyl	1.503	1.380	8.2	94	-0.01
15		Acenaphthylene	9.019	8.456	6.2	96	0.00
16		Acenaphthene	1.300	1.236	4.9	96	-0.01
17		Fluorene	1.768	1.609	9.0	93	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	93	-0.02
19		4-Bromophenyl-phenylether	0.213	0.186	12.7	83	-0.02
20		Hexachlorobenzene	0.885	0.831	6.1	91	-0.02
21		Pentachlorophenol	0.086	0.060	30.2#	68	0.00
22		Phenanthrene	1.281	1.154	9.9	86	0.00
23		Anthracene	1.172	1.074	8.4	87	0.00
24		Fluoranthene	1.431	1.313	8.2	87	-0.01
25	I	Chrysene-d12	1.000	1.000	0.0	90	-0.01
26		Pyrene	1.192	1.134	4.9	87	-0.01
27	S	Terphenyl-d14	0.833	0.767	7.9	83	-0.01
28		Benzo(a)anthracene	1.249	1.195	4.3	88	-0.01
29		Chrysene	1.297	1.179	9.1	83	-0.01
30		Bis(2-ethylhexyl)phthalate	0.398	0.421	-5.8	104	-0.01
31		Indeno(1,2,3-cd)pyrene	1.243	1.245	-0.2	92	-0.03
32	I	Perylene-d12	1.000	1.000	0.0	87	-0.02
33		Benzo(b)fluoranthene	1.103	1.203	-9.1	96	-0.01
34		Benzo(k)fluoranthene	1.232	1.242	-0.8	88	-0.01
35	C	Benzo(a)pyrene	1.019	1.090	-7.0	93	-0.02
36		Dibenzo(a,h)anthracene	1.046	1.054	-0.8	89	-0.03
37		Benzo(g,h,i)perylene	1.060	1.066	-0.6	88	-0.03

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