Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089997.D
Acq On : 30 Jun 2015 1:09
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 30 07:17:49 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	89	0.00
2		1,4-Dioxane	0.686	0.669	2.5	90	0.00
3		n-Nitrosodimethylamine	0.930	0.845	9.1	84	0.00
4	S	2-Fluorophenol	1.149	0.902	21.5	70	0.00
5	S	Phenol-d6	1.608	1.254	22.0	70	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
7	S	Nitrobenzene-d5	0.397	0.344	13.4	78	0.00
8		Nitrobenzene	0.421	0.368	12.6	79	0.00
9		Naphthalene	1.130	1.086	3.9	86	0.00
10		Hexachlorobutadiene	0.178	0.181	-1.7	91	0.00
11		2-Methylnaphthalene	0.755	0.717	5.0	85	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
13	S	2,4,6-Tribromophenol	0.147	0.088	40.1#	54	0.00
14	S	2-Fluorobiphenyl	1.503	1.462	2.7	87	0.00
15		Acenaphthylene	9.019	8.535	5.4	85	0.00
16		Acenaphthene	1.300	1.263	2.8	86	0.00
17		Fluorene	1.768	1.695	4.1	86	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
19		4-Bromophenyl-phenylether	0.213	0.206	3.3	85	0.00
20		Hexachlorobenzene	0.885	0.887	-0.2	90	0.00
21		Pentachlorophenol	0.086	0.047	45.3#	50	0.00
22		Phenanthrene	1.281	1.233	3.7	86	0.00
23		Anthracene	1.172	1.128	3.8	85	0.00
24		Fluoranthene	1.431	1.363	4.8	84	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	88	0.00
26		Pyrene	1.192	1.132	5.0	85	0.00
27	S	Terphenyl-d14	0.833	0.794	4.7	84	0.00
28		Benzo(a)anthracene	1.249	1.162	7.0	84	0.00
29		Chrysene	1.297	1.251	3.5	86	0.00
30		Bis(2-ethylhexyl)phthalate	0.398	0.172	56.8#	41#	0.00
31		Indeno(1,2,3-cd)pyrene	1.243	1.050	15.5	75	0.00
32	I	Perylene-d12	1.000	1.000	0.0	84	0.00
33		Benzo(b)fluoranthene	1.103	1.066	3.4	81	0.00
34		Benzo(k)fluoranthene	1.232	1.191	3.3	81	0.00
35	C	Benzo(a)pyrene	1.019	0.930	8.7	76	0.00
36		Dibenzo(a,h)anthracene	1.046	0.927	11.4	75	0.00
37		Benzo(g,h,i)perylene	1.060	0.955	9.9	76	0.00

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