Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071015\
Data File : BE090083.D
Acq On : 10 Jul 2015 15:04
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 10 23:13:55 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	120	-0.01
2		1,4-Dioxane	0.686	0.542	21.0	99	-0.01
3		n-Nitrosodimethylamine	0.930	0.731	21.4	98	-0.01
4	S	2-Fluorophenol	1.149	1.282	-11.6	135	0.00
5	S	Phenol-d6	1.608	2.349	-46.1#	179#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	127	-0.01
7	S	Nitrobenzene-d5	0.397	0.414	-4.3	135	0.00
8		Nitrobenzene	0.421	0.339	19.5	106	0.00
9		Naphthalene	1.130	0.985	12.8	113	-0.01
10		Hexachlorobutadiene	0.178	0.160	10.1	117	0.00
11		2-Methylnaphthalene	0.755	0.626	17.1	107	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
13	S	2,4,6-Tribromophenol	0.147	0.202	-37.4#	163#	0.00
14	S	2-Fluorobiphenyl	1.503	1.846	-22.8	146	-0.01
15		Acenaphthylene	9.019	7.740	14.2	102	0.00
16		Acenaphthene	1.300	1.126	13.4	102	-0.01
17		Fluorene	1.768	1.461	17.4	98	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	101	-0.02
19		4-Bromophenyl-phenylether	0.213	0.180	15.5	88	-0.02
20		Hexachlorobenzene	0.885	0.749	15.4	89	-0.02
21		Pentachlorophenol	0.086	0.065	24.4	81	0.00
22		Phenanthrene	1.281	1.069	16.5	87	-0.02
23		Anthracene	1.172	0.973	17.0	86	0.00
24		Fluoranthene	1.431	1.138	20.5	82	-0.02
25	I	Chrysene-d12	1.000	1.000	0.0	88	-0.01
26		Pyrene	1.192	1.092	8.4	82	-0.01
27	S	Terphenyl-d14	0.833	1.098	-31.8#	116	-0.01
28		Benzo(a)anthracene	1.249	1.072	14.2	77	-0.01
29		Chrysene	1.297	1.084	16.4	75	-0.01
30		Bis(2-ethylhexyl)phthalate	0.398	0.366	8.0	88	-0.02
31		Indeno(1,2,3-cd)pyrene	1.243	1.044	16.0	75	-0.03
32	I	Perylene-d12	1.000	1.000	0.0	82	-0.02
33		Benzo(b)fluoranthene	1.103	1.020	7.5	77	-0.02
34		Benzo(k)fluoranthene	1.232	1.154	6.3	77	-0.02
35	C	Benzo(a)pyrene	1.019	0.940	7.8	76	-0.02
36		Dibenzo(a,h)anthracene	1.046	0.919	12.1	73	-0.03
37		Benzo(g,h,i)perylene	1.060	0.935	11.8	73	-0.04

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