Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE091615\
Data File : BE090842.D
Acq On : 16 Sep 2015 11:32
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 17 00:54:42 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	110	0.00
2		1,4-Dioxane	0.814	0.643	21.0	90	0.00
3		n-Nitrosodimethylamine	1.033	0.804	22.2	91	0.00
4	S	2-Fluorophenol	1.039	0.756	27.2#	85	0.00
5	S	Phenol-d6	1.392	1.258	9.6	105	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
7	S	Nitrobenzene-d5	0.330	0.284	13.9	108	0.00
8		Nitrobenzene	0.365	0.321	12.1	115	0.00
9		Naphthalene	1.091	0.954	12.6	110	-0.01
10		Hexachlorobutadiene	0.171	0.144	15.8	103	0.00
11		2-Methylnaphthalene	0.590	0.532	9.8	117	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	138	0.00
13	S	2,4,6-Tribromophenol	0.143	0.089	37.8#	102	0.00
14	S	2-Fluorobiphenyl	1.385	1.131	18.3	115	-0.01
15		Acenaphthylene	7.992	6.744	15.6	122	0.00
16		Acenaphthene	1.335	1.154	13.6	118	0.00
17		Fluorene	1.666	1.419	14.8	119	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	133	0.00
19		4-Bromophenyl-phenylether	0.170	0.135	20.6	109	-0.02
20		Hexachlorobenzene	0.928	0.769	17.1	111	0.00
21		Pentachlorophenol	0.054	0.036	33.3#	101	0.00
22		Phenanthrene	1.237	1.100	11.1	118	-0.02
23		Anthracene	1.035	0.911	12.0	122	0.00
24		Fluoranthene	1.304	1.135	13.0	124	-0.01
25	I	Chrysene-d12	1.000	1.000	0.0	145	0.00
26		Pyrene	1.035	0.999	3.5	148	-0.01
27	S	Terphenyl-d14	0.553	0.478	13.6	131	-0.01
28		Benzo(a)anthracene	1.156	1.030	10.9	136	0.00
29		Chrysene	1.279	1.131	11.6	130	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.238	31.0#	136	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.214	-2.8	161#	-0.02
32	I	Perylene-d12	1.000	1.000	0.0	160#	-0.01
33		Benzo(b)fluoranthene	1.084	0.966	10.9	147	-0.01
34		Benzo(k)fluoranthene	1.263	1.165	7.8	150#	-0.01
35	C	Benzo(a)pyrene	0.970	0.911	6.1	161#	-0.01
36		Dibenzo(a,h)anthracene	1.020	0.971	4.8	159#	-0.02
37		Benzo(g,h,i)perylene	1.116	0.959	14.1	146	-0.03

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