Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090815\
Data File : BE090755.D
Acq On : 8 Sep 2015 19:15
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 08 23:08:18 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 08 23:05:35 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	133	0.00
2		1,4-Dioxane	0.814	0.703	13.6	118	0.00
3		n-Nitrosodimethylamine	1.033	0.906	12.3	123	0.00
4	S	2-Fluorophenol	1.039	0.997	4.0	136	0.00
5	S	Phenol-d6	1.392	1.581	-13.6	159#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	148	0.00
7	S	Nitrobenzene-d5	0.330	0.339	-2.7	152#	0.00
8		Nitrobenzene	0.365	0.376	-3.0	158#	0.00
9		Naphthalene	1.091	1.101	-0.9	150	0.00
10		Hexachlorobutadiene	0.171	0.164	4.1	137	0.00
11		2-Methylnaphthalene	0.590	0.664	-12.5	171#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	172#	0.00
13	S	2,4,6-Tribromophenol	0.143	0.142	0.7	202#	0.00
14	S	2-Fluorobiphenyl	1.385	1.341	3.2	169#	0.00
15		Acenaphthylene	7.992	8.245	-3.2	185#	0.00
16		Acenaphthene	1.335	1.332	0.2	170#	0.00
17		Fluorene	1.666	1.686	-1.2	176#	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	164#	0.00
19		4-Bromophenyl-phenylether	0.170	0.176	-3.5	176#	-0.02
20		Hexachlorobenzene	0.928	0.895	3.6	160#	0.00
21		Pentachlorophenol	0.054	0.051	5.6	178#	0.00
22		Phenanthrene	1.237	1.274	-3.0	169#	-0.02
23		Anthracene	1.035	1.151	-11.2	190#	0.00
24		Fluoranthene	1.304	1.345	-3.1	182#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	157#	0.00
26		Pyrene	1.035	1.207	-16.6	193#	0.00
27	S	Terphenyl-d14	0.553	0.619	-11.9	184#	0.00
28		Benzo(a)anthracene	1.156	1.203	-4.1	173#	0.00
29		Chrysene	1.279	1.337	-4.5	166#	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.320	7.2	198#	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.339	-13.4	193#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	167#	0.00
33		Benzo(b)fluoranthene	1.084	1.157	-6.7	183#	0.00
34		Benzo(k)fluoranthene	1.263	1.310	-3.7	175#	0.00
35	C	Benzo(a)pyrene	0.970	1.049	-8.1	193#	0.00
36		Dibenzo(a,h)anthracene	1.020	1.124	-10.2	191#	0.00
37		Benzo(g,h,i)perylene	1.116	1.141	-2.2	180#	-0.01

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