Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081315\
Data File : BE090499.D
Acq On : 13 Aug 2015 13:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 13 15:11:26 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 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	130	-0.01
2		1,4-Dioxane	0.687	0.709	-3.2	125	0.00
3		n-Nitrosodimethylamine	0.773	0.865	-11.9	146	0.00
4	S	2-Fluorophenol	0.750	1.496	-99.5#	278#	0.00
5	S	Phenol-d6	1.241	2.205	-77.7#	249#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	145	0.00
7	S	Nitrobenzene-d5	0.331	0.403	-21.8	186#	-0.01
8		Nitrobenzene	0.349	0.422	-20.9	196#	-0.01
9		Naphthalene	1.104	1.132	-2.5	150	0.00
10		Hexachlorobutadiene	0.191	0.166	13.1	125	0.00
11		2-Methylnaphthalene	0.640	0.742	-15.9	176#	-0.01
12	I	Acenaphthene-d10	1.000	1.000	0.0	144	0.00
13	S	2,4,6-Tribromophenol	0.095	0.207	-117.9#	402#	-0.02
14	S	2-Fluorobiphenyl	1.375	1.503	-9.3	155#	0.00
15		Acenaphthylene	8.098	9.116	-12.6	163#	0.00
16		Acenaphthene	1.259	1.372	-9.0	156#	0.00
17		Fluorene	1.635	1.683	-2.9	146	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	134	0.00
19		4-Bromophenyl-phenylether	0.194	0.206	-6.2	146	0.00
20		Hexachlorobenzene	1.005	0.904	10.0	120	0.00
21		Pentachlorophenol	0.044	0.081	-84.1#	401#	0.00
22		Phenanthrene	1.251	1.303	-4.2	141	0.00
23		Anthracene	1.028	1.209	-17.6	162#	-0.02
24		Fluoranthene	1.231	1.420	-15.4	161#	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	118	0.00
26		Pyrene	1.084	1.449	-33.7#	164#	0.00
27	S	Terphenyl-d14	0.737	0.925	-25.5#	150	0.00
28		Benzo(a)anthracene	1.108	1.297	-17.1	154#	0.00
29		Chrysene	1.339	1.337	0.1	115	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.746	-139.9#	395#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.282	-36.1#	195#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	128	0.00
33		Benzo(b)fluoranthene	0.918	1.202	-30.9#	178#	-0.01
34		Benzo(k)fluoranthene	1.392	1.350	3.0	121	0.00
35	C	Benzo(a)pyrene	0.850	1.112	-30.8#	178#	0.00
36		Dibenzo(a,h)anthracene	0.766	1.076	-40.5#	208#	-0.01
37		Benzo(g,h,i)perylene	0.923	1.192	-29.1#	171#	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1