Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081115\
Data File : BE090443.D
Acq On : 12 Aug 2015 00:13
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 12 05:46:15 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	120	0.00
2		1,4-Dioxane	0.687	0.641	6.7	104	0.00
3		n-Nitrosodimethylamine	0.773	0.785	-1.6	122	0.00
4	S	2-Fluorophenol	0.750	1.052	-40.3#	180#	0.00
5	S	Phenol-d6	1.241	1.604	-29.3#	167#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	126	0.00
7	S	Nitrobenzene-d5	0.331	0.348	-5.1	139	0.00
8		Nitrobenzene	0.349	0.364	-4.3	146	0.00
9		Naphthalene	1.104	1.124	-1.8	129	0.00
10		Hexachlorobutadiene	0.191	0.192	-0.5	125	0.00
11		2-Methylnaphthalene	0.640	0.701	-9.5	144	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	134	0.00
13	S	2,4,6-Tribromophenol	0.095	0.130	-36.8#	236#	0.00
14	S	2-Fluorobiphenyl	1.375	1.445	-5.1	139	0.00
15		Acenaphthylene	8.098	8.517	-5.2	142	0.00
16		Acenaphthene	1.259	1.275	-1.3	136	0.00
17		Fluorene	1.635	1.683	-2.9	137	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	129	0.00
19		4-Bromophenyl-phenylether	0.194	0.200	-3.1	136	0.00
20		Hexachlorobenzene	1.005	0.942	6.3	120	0.00
21		Pentachlorophenol	0.044	0.042	4.5	201#	0.00
22		Phenanthrene	1.251	1.292	-3.3	134	0.00
23		Anthracene	1.028	1.147	-11.6	147	-0.02
24		Fluoranthene	1.231	1.316	-6.9	143	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	127	0.00
26		Pyrene	1.084	1.189	-9.7	145	0.00
27	S	Terphenyl-d14	0.737	0.823	-11.7	144	-0.01
28		Benzo(a)anthracene	1.108	1.175	-6.0	150#	0.00
29		Chrysene	1.339	1.348	-0.7	125	0.00
30		Bis(2-ethylhexyl)phthalate	0.311	0.322	-3.5	184#	0.00
31		Indeno(1,2,3-cd)pyrene	0.942	1.014	-7.6	166#	0.00
32	I	Perylene-d12	1.000	1.000	0.0	132	0.00
33		Benzo(b)fluoranthene	0.918	1.050	-14.4	160#	0.00
34		Benzo(k)fluoranthene	1.392	1.318	5.3	121	0.00
35	C	Benzo(a)pyrene	0.850	0.925	-8.8	153#	0.00
36		Dibenzo(a,h)anthracene	0.766	0.868	-13.3	173#	0.00
37		Benzo(g,h,i)perylene	0.923	0.981	-6.3	145	0.00

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