Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE082615\
Data File : BE090645.D
Acq On : 26 Aug 2015 21:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 26 23:51:16 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 26 23:44:39 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	102	0.00
2		1,4-Dioxane	0.658	0.731	-11.1	107	0.00
3		n-Nitrosodimethylamine	0.863	0.872	-1.0	104	0.00
4	S	2-Fluorophenol	1.446	1.029	28.8#	72	0.00
5	S	Phenol-d6	2.125	1.467	31.0#	73	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
7	S	Nitrobenzene-d5	0.398	0.364	8.5	90	0.00
8		Nitrobenzene	0.413	0.365	11.6	88	0.00
9		Naphthalene	1.136	1.101	3.1	91	0.00
10		Hexachlorobutadiene	0.167	0.174	-4.2	97	0.00
11		2-Methylnaphthalene	0.740	0.678	8.4	87	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
13	S	2,4,6-Tribromophenol	0.192	0.156	18.8	91	0.00
14	S	2-Fluorobiphenyl	1.493	1.341	10.2	92	0.00
15		Acenaphthylene	9.132	8.319	8.9	95	0.00
16		Acenaphthene	1.374	1.338	2.6	100	0.00
17		Fluorene	1.684	1.761	-4.6	107	0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.00
19		4-Bromophenyl-phenylether	0.204	0.180	11.8	106	0.00
20		Hexachlorobenzene	0.906	0.904	0.2	118	-0.02
21		Pentachlorophenol	0.079	0.057	27.8#	113	0.00
22		Phenanthrene	1.312	1.283	2.2	117	0.00
23		Anthracene	1.208	1.137	5.9	114	0.00
24		Fluoranthene	1.396	1.367	2.1	118	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	138	0.00
26		Pyrene	1.430	1.255	12.2	124	0.00
27	S	Terphenyl-d14	0.914	0.793	13.2	122	0.00
28		Benzo(a)anthracene	1.316	1.182	10.2	130	0.00
29		Chrysene	1.344	1.371	-2.0	142	0.00
30		Bis(2-ethylhexyl)phthalate	0.725	0.386	46.8#	89	0.00
31		Indeno(1,2,3-cd)pyrene	1.266	1.145	9.6	135	0.00
32	I	Perylene-d12	1.000	1.000	0.0	138	0.00
33		Benzo(b)fluoranthene	1.184	1.095	7.5	131	0.00
34		Benzo(k)fluoranthene	1.342	1.369	-2.0	140	0.00
35	C	Benzo(a)pyrene	1.108	1.024	7.6	132	0.00
36		Dibenzo(a,h)anthracene	1.069	0.958	10.4	130	0.00
37		Benzo(g,h,i)perylene	1.169	1.059	9.4	127	0.00

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