Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080816\
Data File : BE092232.D
Acq On : 9 Aug 2016 1:14
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 09 02:44:37 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-PAH-BE080216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 08 18:55:30 2016
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	87	0.00
2		1,4-Dioxane	0.680	0.670	1.5	83	0.00
3		n-Nitrosodimethylamine	0.942	0.761	19.2	87	0.00
4	S	2-Fluorophenol	1.284	1.119	12.9	86	0.01
5	S	Phenol-d6	1.711	1.490	12.9	82	0.01
6	I	Naphthalene-d8	1.000	1.000	0.0	93	-0.01
7	S	Nitrobenzene-d5	0.395	0.322	18.5	80	0.00
8		Naphthalene	1.159	1.113	4.0	92	0.00
9		2-Methylnaphthalene	0.743	0.700	5.8	94	0.00
10	I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.02
11	S	2,4,6-Tribromophenol	0.162	0.137	15.4	95	0.00
12	S	2-Fluorobiphenyl	1.584	1.531	3.3	95	0.00
13		Acenaphthylene	8.821	8.367	5.1	95	-0.02
14		Acenaphthene	1.406	1.365	2.9	94	0.00
15		Fluorene	1.783	1.663	6.7	93	-0.01
16	I	Phenanthrene-d10	1.000	1.000	0.0	95	-0.02
17		Hexachlorobenzene	0.245	0.229	6.5	88	0.00
18		Pentachlorophenol	0.068	0.053	22.1	91	0.00
19		Phenanthrene	1.460	1.412	3.3	96	0.00
20		Anthracene	1.331	1.218	8.5	92	0.00
21		Fluoranthene	6.116	5.627	8.0	91	0.00
22	I	Chrysene-d12	1.000	1.000	0.0	83	0.00
23		Pyrene	5.228	5.227	0.0	90	0.00
24	S	Terphenyl-d14	0.718	0.753	-4.9	95	-0.02
25		Benzo(a)anthracene	5.496	5.501	-0.1	89	0.00
26		Chrysene	5.505	5.378	2.3	86	0.00
27		Indeno(1,2,3-cd)pyrene	5.978	5.455	8.7	82	0.00
28	I	Perylene-d12	1.000	1.000	0.0	82	0.00
29		Benzo(b)fluoranthene	1.368	1.285	6.1	76	0.00
30		Benzo(k)fluoranthene	1.369	1.328	3.0	92	0.00
31	C	Benzo(a)pyrene	1.257	1.235	1.8	84	0.00
32		Dibenzo(a,h)anthracene	1.200	1.126	6.2	82	0.00
33		Benzo(g,h,i)perylene	5.122	4.812	6.1	81	0.00

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