

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111615\
 Data File : BE091186.D
 Acq On : 16 Nov 2015 17:59
 Operator : UM/NP
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTD0.469

Quant Time: Nov 17 10:35:08 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM02.2-EPA-SIM-BE102715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 16 04:54:02 2015
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	115	0.00
2 SURR	1,4-Dioxane-d8	0.288	0.251	12.8	93	0.05
3	1,4-Dioxane	0.325	0.270	16.9	93	0.05
4 I	Naphthalene-d8	1.000	1.000	0.0	119	0.00
5	Naphthalene	0.986	0.948	3.9	111	0.00
6 SURR	2-Methylnaphthalene-d10	0.546	0.510	6.6	111	0.00
7	2-Methylnaphthalene	0.643	0.582	9.5	107	0.00
8 I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
9	Acenaphthylene	1.998	1.911	4.4	107	0.00
10 C	Acenaphthene	1.368	1.346	1.6	111	0.00
11	Fluorene	1.757	1.728	1.7	106	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	106	0.01
13	Pentachlorophenol	0.038	0.042	-10.5	145	0.00
14	Phenanthrene	4.619	4.513	2.3	102	0.00
15	Anthracene	4.200	4.032	4.0	105	0.00
16 SURR	Fluoranthene-d10	1.127	1.100	2.4	110	0.00
17 C	Fluoranthene	5.406	5.352	1.0	110	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	115	0.00
19	Pyrene	4.544	4.272	6.0	107	0.00
20	Benzo(a)anthracene	1.027	0.962	6.3	111	0.00
21	Chrysene	1.183	1.167	1.4	112	0.00
22 I	Perylene-d12	1.000	1.000	0.0	121	0.00
23	Benzo(b)fluoranthene	1.249	1.214	2.8	115	0.00
24	Benzo(k)fluoranthene	1.281	1.243	3.0	112	0.00
25 C	Benzo(a)pyrene	1.150	1.224	-6.4	126	0.00
26	Indeno(1,2,3-cd)pyrene	4.847	4.623	4.6	111	0.01
27	Dibenzo(a,h)anthracene	1.155	1.145	0.9	119	0.01
28	Benzo(q,h,i)perylene	5.083	4.868	4.2	114	0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0