

Data Path : Z:\HPCHEM1\BNA E\Data\BE111615\
 Data File : BE091176.D
 Acq On : 16 Nov 2015 10:07
 Operator : UM/NP
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTD0.468

Quant Time: Nov 17 10:31:18 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	125	0.02
2 SURR	1,4-Dioxane-d8	0.288	0.238	17.4	97	0.06
3	1,4-Dioxane	0.325	0.271	16.6	102	0.06
4 I	Naphthalene-d8	1.000	1.000	0.0	128	0.02
5	Naphthalene	0.986	0.946	4.1	119	0.02
6 SURR	2-Methylnaphthalene-d10	0.546	0.509	6.8	119	0.02
7	2-Methylnaphthalene	0.643	0.578	10.1	114	0.01
8 I	Acenaphthene-d10	1.000	1.000	0.0	123	0.00
9	Acenaphthylene	1.998	1.884	5.7	112	0.00
10 C	Acenaphthene	1.368	1.341	2.0	117	0.00
11	Fluorene	1.757	1.655	5.8	108	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	108	0.01
13	Pentachlorophenol	0.038	0.034	10.5	120	0.01
14	Phenanthrene	4.619	4.414	4.4	101	0.01
15	Anthracene	4.200	4.065	3.2	107	0.01
16 SURR	Fluoranthene-d10	1.127	1.098	2.6	112	0.00
17 C	Fluoranthene	5.406	5.390	0.3	112	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	119	0.00
19	Pyrene	4.544	4.198	7.6	109	0.01
20	Benzo(a)anthracene	1.027	0.961	6.4	114	0.00
21	Chrysene	1.183	1.167	1.4	116	0.00
22 I	Perylene-d12	1.000	1.000	0.0	127	0.00
23	Benzo(b)fluoranthene	1.249	1.172	6.2	117	0.00
24	Benzo(k)fluoranthene	1.281	1.253	2.2	119	0.00
25 C	Benzo(a)pyrene	1.150	1.202	-4.5	131	0.01
26	Indeno(1,2,3-cd)pyrene	4.847	4.672	3.6	119	0.02
27	Dibenzo(a,h)anthracene	1.155	1.134	1.8	124	0.01
28	Benzo(q,h,i)perylene	5.083	4.893	3.7	121	0.02

(#) = Out of Range

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