

Data Path : Z:\HPCHEM1\BNA E\Data\BE112115\
 Data File : BE091258.D
 Acq On : 21 Nov 2015 00:05
 Operator : UM/NP
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTD0.479

Quant Time: Nov 21 04:53:33 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 : Sat Nov 21 02:52:22 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	68	0.00
2 SURR	1,4-Dioxane-d8	0.288	0.305	-5.9	67	0.02
3	1,4-Dioxane	0.325	0.344	-5.8	70	0.02
4 I	Naphthalene-d8	1.000	1.000	0.0	61	0.02
5	Naphthalene	0.986	1.015	-2.9	61	0.02
6 SURR	2-Methylnaphthalene-d10	0.546	0.472	13.6	52	0.02
7	2-Methylnaphthalene	0.643	0.530	17.6	49	0.02
8 I	Acenaphthene-d10	1.000	1.000	0.0	44	0.00
9	Acenaphthylene	1.998	2.133	-6.8	46	0.00
10 C	Acenaphthene	1.368	1.523	-11.3	48	0.00
11	Fluorene	1.757	1.942	-10.5	45	0.00
12 I	Phenanthrene-d10	1.000	1.000	0.0	43	0.01
13	Pentachlorophenol	0.038	0.046	-21.1#	66	0.00
14	Phenanthrene	4.619	4.645	-0.6	43	0.00
15	Anthracene	4.200	4.059	3.4	43	0.00
16 SURR	Fluoranthene-d10	1.127	1.017	9.8	42	0.00
17 C	Fluoranthene	5.406	5.468	-1.1	46	0.00
18 I	Chrysene-d12	1.000	1.000	0.0	46	0.00
19	Pyrene	4.544	4.149	8.7	42	0.00
20	Benzo(a)anthracene	1.027	0.881	14.2	41	0.00
21	Chrysene	1.183	1.242	-5.0	48	0.00
22 I	Perylene-d12	1.000	1.000	0.0	42	0.00
23	Benzo(b)fluoranthene	1.249	1.342	-7.4	44	0.00
24	Benzo(k)fluoranthene	1.281	1.518	-18.5	47	0.00
25 C	Benzo(a)pyrene	1.150	1.248	-8.5	44	0.00
26	Indeno(1,2,3-cd)pyrene	4.847	5.353	-10.4	44	0.00
27	Dibenzo(a,h)anthracene	1.155	1.211	-4.8	43	0.00
28	Benzo(q,h,i)perylene	5.083	5.568	-9.5	45	0.00

(#) = Out of Range

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