

Data Path : Z:\HPCHEM1\BNA E\Data\BE042716\
 Data File : BE091722.D
 Acq On : 27 Apr 2016 12:34
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 28 03:18:22 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 26 16:09:46 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
2	1,4-Dioxane	0.752	0.667	11.3	85	0.00
3	n-Nitrosodimethylamine	0.908	0.807	11.1	88	0.02
4 S	2-Fluorophenol	1.228	1.127	8.2	92	0.00
5 S	Phenol-d6	1.629	1.408	13.6	89	0.00
6	bis(2-Chloroethyl)ether	1.433	1.371	4.3	92	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
8 S	Nitrobenzene-d5	0.323	0.283	12.4	90	0.00
9	Nitrobenzene	0.338	0.288	14.8	89	0.00
10	Naphthalene	1.010	1.024	-1.4	96	0.00
11	Hexachlorobutadiene	0.149	0.151	-1.3	99	0.00
12	2-Methylnaphthalene	0.649	0.630	2.9	93	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	93	-0.01
14 S	2,4,6-Tribromophenol	0.157	0.127	19.1	82	0.00
15 S	2-Fluorobiphenyl	1.414	1.417	-0.2	94	0.00
16	Acenaphthylene	1.923	1.894	1.5	104	0.00
17	Acenaphthene	1.247	1.245	0.2	93	0.00
18	Fluorene	1.550	1.522	1.8	91	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
20	4-Bromophenyl-phenylether	0.180	0.174	3.3	88	0.00
21	Hexachlorobenzene	0.180	0.179	0.6	92	0.00
22	Pentachlorophenol	0.085	0.067	21.2	86	0.00
23	Phenanthrene	1.136	1.155	-1.7	92	0.00
24	Anthracene	1.025	0.965	5.9	89	0.00
25	Fluoranthene	1.188	1.157	2.6	89	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	86	0.00
27	Benzydine	0.286	0.292	-2.1	108	0.00
28	Pvrene	1.130	1.154	-2.1	91	0.00
29 S	Terphenyl-d14	0.780	0.795	-1.9	89	0.00
30	Benzo(a)anthracene	1.221	1.215	0.5	87	0.00
31	3,3'-Dichlorobenzidine	0.623	0.545	12.5	87	-0.02
32	Chrysene	1.269	1.517	-19.5	97	0.00
33	Bis(2-ethylhexyl)phthalate	0.370	0.323	12.7	97	-0.02
34	Indeno(1,2,3-cd)pyrene	1.130	1.247	-10.4	93	0.00
35 I	Perylene-d12	1.000	1.000	0.0	93	0.00
36	Benzo(b)fluoranthene	1.172	1.204	-2.7	99	0.00
37	Benzo(k)fluoranthene	1.191	1.100	7.6	90	0.00
38 C	Benzo(a)pyrene	1.103	1.047	5.1	94	0.00
39	Dibenzo(a,h)anthracene	1.045	0.999	4.4	92	-0.01
40	Benzo(a,h,i)perylene	1.062	1.045	1.6	95	-0.01

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(#) = Out of Range

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