

Data Path : Z:\HPCHEM1\BNA_E\Data\BE111617\
 Data File : BE094818.D
 Acq On : 15 Nov 2017 17:09
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 16 02:01:26 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 22:40:47 2017
 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	66	0.01
2	1,4-Dioxane	0.677	0.767	-13.3	65	-0.03
3	n-Nitrosodimethylamine	0.648	0.578	10.8	59	0.00
4 S	2-Fluorophenol	1.370	1.179	13.9	57	0.02
5 S	Phenol-d6	2.069	2.015	2.6	66	0.06
6	bis(2-Chloroethyl)ether	1.600	1.566	2.1	66	0.02
7 I	Naphthalene-d8	1.000	1.000	0.0	69	0.02
8 S	Nitrobenzene-d5	0.319	0.286	10.3	64	0.05
9	Naphthalene	4.527	4.051	10.5	60	0.03
10	Hexachlorobutadiene	0.164	0.169	-3.0	69	-0.02
11 SURR	2-Methylnaphthalene-d10	0.642	0.577	10.1	60	0.03
12	2-Methylnaphthalene	0.768	0.682	11.2	60	0.03
13 I	Acenaphthene-d10	1.000	1.000	0.0	73	0.02
14 S	2,4,6-Tribromophenol	0.110	0.142	-29.1#	97	0.03
15 S	2-Fluorobiphenyl	1.410	1.222	13.3	62	0.02
16	Acenaphthylene	2.112	1.861	11.9	65	0.00
17	Acenaphthene	1.356	1.222	9.9	65	0.00
18	Fluorene	1.738	1.673	3.7	70	0.02
19 I	Phenanthrene-d10	1.000	1.000	0.0	94	0.03
20	4-Bromophenyl-phenylether	0.218	0.163	25.2#	71	0.03
21	Hexachlorobenzene	0.191	0.234	-22.5	100	0.00
22	Pentachlorophenol	0.024	0.026	-8.3	100	0.03
23	Phenanthrene	1.141	1.329	-16.5	108	0.00
24	Anthracene	1.176	1.325	-12.7	103	0.00
25 SURR	Fluoranthene-d10	4.676	5.229	-11.8	91	0.01
26	Fluoranthene	1.420	1.566	-10.3	90	0.01
27 I	Chrysene-d12	1.000	1.000	0.0	99	0.01
28	Pyrene	1.399	1.271	9.1	91	0.01
29 S	Terphenyl-d14	0.754	0.686	9.0	91	0.00
30	Benzo(a)anthracene	1.302	1.211	7.0	92	0.01
31	Chrysene	1.322	1.313	0.7	101	0.01
32	Bis(2-ethylhexyl)phthalate	3.193	2.873	10.0	97	-0.01
33	Indeno(1,2,3-cd)pyrene	1.222	1.197	2.0	98	0.07
34 I	Perylene-d12	1.000	1.000	0.0	98	0.03
35	Benzo(b)fluoranthene	1.294	1.173	9.4	92	0.02
36	Benzo(k)fluoranthene	1.379	1.288	6.6	93	0.03
37 C	Benzo(a)pyrene	1.255	1.202	4.2	94	0.04
38	Dibenzo(a,h)anthracene	1.127	1.110	1.5	99	0.05
39	Benzo(g,h,i)perylene	1.055	1.102	-4.5	104	0.09

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