

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120517\
 Data File : BE094931.D
 Acq On : 6 Dec 2017 00:59
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 06 06:54:43 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 01 13:33:28 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	79	0.00
2	1,4-Dioxane	0.400	0.372	7.0	69	0.00
3	n-Nitrosodimethylamine	0.400	0.430	-7.5	86	0.00
4 S	2-Fluorophenol	0.400	0.443	-10.7	86	0.00
5 S	Phenol-d6	0.400	0.439	-9.7	86	0.00
6	bis(2-Chloroethyl)ether	0.400	0.383	4.3	74	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	80	0.00
8 S	Nitrobenzene-d5	0.400	0.555	-38.8#	112	0.00
9	Naphthalene	0.400	0.386	3.5	77	0.00
10	Hexachlorobutadiene	0.400	0.408	-2.0	81	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	81	0.00
12	2-Methylnaphthalene	0.400	0.401	-0.3	81	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	88	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.676	-69.0#	166	0.00
15 S	2-Fluorobiphenyl	0.400	0.376	6.0	81	0.00
16	Acenaphthylene	0.400	0.383	4.3	86	0.00
17	Acenaphthene	0.400	0.385	3.8	86	0.00
18	Fluorene	0.400	0.385	3.8	83	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.400	0.408	-2.0	89	0.00
21	Hexachlorobenzene	0.400	0.404	-1.0	86	0.00
22	Pentachlorophenol	0.400	0.634	-58.5#	160	0.00
23	Phenanthrene	0.400	0.382	4.5	82	0.00
24	Anthracene	0.400	0.380	5.0	85	0.00
25 SURR	Fluoranthene-d10	0.400	0.390	2.5	85	0.00
26	Fluoranthene	0.400	0.377	5.8	82	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	73	0.00
28	Pvrene	0.400	0.446	-11.5	82	-0.02
29 S	Terphenyl-d14	0.400	0.444	-11.0	82	0.00
30	Benzo(a)anthracene	0.400	0.439	-9.7	82	-0.02
31	Chrysene	0.400	0.375	6.3	69	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.646	-61.5#	138	-0.02
33	Indeno(1,2,3-cd)pyrene	0.400	0.429	-7.2	80	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	82	0.00
35	Benzo(b)fluoranthene	0.400	0.379	5.3	80	0.00
36	Benzo(k)fluoranthene	0.400	0.360	10.0	76	0.00
37 C	Benzo(a)pyrene	0.400	0.386	3.5	82	0.00
38	Dibenzo(a,h)anthracene	0.400	0.382	4.5	81	-0.01
39	Benzo(a,h,i)perylene	0.400	0.380	5.0	80	-0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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