

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA E\Data\BE120517\
 Data File : BE094948.D
 Acq On : 6 Dec 2017 12:43
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 06 16:48:15 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	84	-0.01
2	1,4-Dioxane	0.400	0.390	2.5	77	0.00
3	n-Nitrosodimethylamine	0.400	0.366	8.5	79	-0.01
4 S	2-Fluorophenol	0.400	0.408	-2.0	85	0.00
5 S	Phenol-d6	0.400	0.405	-1.3	85	0.00
6	bis(2-Chloroethyl)ether	0.400	0.368	8.0	76	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	81	0.00
8 S	Nitrobenzene-d5	0.400	0.524	-31.0#	107	-0.02
9	Naphthalene	0.400	0.393	1.8	79	0.00
10	Hexachlorobutadiene	0.400	0.405	-1.3	81	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	82	0.00
12	2-Methylnaphthalene	0.400	0.392	2.0	80	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	85	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.515	-28.7#	114	0.00
15 S	2-Fluorobiphenyl	0.400	0.397	0.8	83	0.00
16	Acenaphthylene	0.400	0.379	5.3	82	0.00
17	Acenaphthene	0.400	0.388	3.0	84	-0.02
18	Fluorene	0.400	0.389	2.8	81	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.400	0.396	1.0	85	0.00
21	Hexachlorobenzene	0.400	0.407	-1.7	85	0.00
22	Pentachlorophenol	0.400	0.427	-6.7	92	-0.01
23	Phenanthrene	0.400	0.383	4.3	81	0.00
24	Anthracene	0.400	0.375	6.3	82	0.00
25 SURR	Fluoranthene-d10	0.400	0.364	9.0	78	0.00
26	Fluoranthene	0.400	0.364	9.0	78	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	76	-0.01
28	Pvrene	0.400	0.396	1.0	76	-0.01
29 S	Terphenyl-d14	0.400	0.402	-0.5	78	-0.01
30	Benzo(a)anthracene	0.400	0.386	3.5	76	-0.01
31	Chrysene	0.400	0.377	5.8	72	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	83	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.378	5.5	73	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	74	-0.01
35	Benzo(b)fluoranthene	0.400	0.391	2.3	74	-0.01
36	Benzo(k)fluoranthene	0.400	0.367	8.3	70	-0.01
37 C	Benzo(a)pyrene	0.400	0.386	3.5	73	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.390	2.5	74	-0.03
39	Benzo(a,h,i)perylene	0.400	0.387	3.3	73	-0.02

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

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