

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE040419\
 Data File : BE099293.D
 Acq On : 4 Apr 2019 16:00
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 05 04:29:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE040219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 02 17:16:42 2019
 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	107	-0.02
2	1,4-Dioxane	0.400	0.445	-11.2	119	-0.01
3	n-Nitrosodimethylamine	0.400	0.346	13.5	95	-0.02
4 S	2-Fluorophenol	0.400	0.381	4.8	97	-0.02
5 S	Phenol-d6	0.400	0.392	2.0	103	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.420	-5.0	112	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	104	-0.02
8 S	Nitrobenzene-d5	0.400	0.422	-5.5	114	-0.02
9	Naphthalene	0.400	0.430	-7.5	111	-0.02
10	Hexachlorobutadiene	0.400	0.423	-5.7	109	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.423	-5.7	112	-0.01
12	2-Methylnaphthalene	0.400	0.417	-4.2	109	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	102	-0.03
14 S	2,4,6-Tribromophenol	0.400	0.342	14.5	97	-0.01
15 S	2-Fluorobiphenyl	0.400	0.440	-10.0	112	-0.01
16	Acenaphthylene	0.400	0.391	2.3	103	-0.01
17	Acenaphthene	0.400	0.414	-3.5	108	-0.01
18	Fluorene	0.400	0.414	-3.5	105	-0.03
19 I	Phenanthrene-d10	0.400	0.400	0.0	97	-0.03
20	4-Bromophenyl-phenylether	0.400	0.423	-5.7	106	-0.03
21	Hexachlorobenzene	0.400	0.442	-10.5	109	-0.03
22	Pentachlorophenol	0.400	0.409	-2.2	114	-0.03
23	Phenanthrene	0.400	0.429	-7.2	103	-0.03
24	Anthracene	0.400	0.397	0.8	97	-0.01
25 SURR	Fluoranthene-d10	0.400	0.402	-0.5	97	-0.02
26	Fluoranthene	0.400	0.412	-3.0	100	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	94	-0.02
28	Pyrene	0.400	0.414	-3.5	99	-0.02
29 S	Terphenyl-d14	0.400	0.421	-5.2	99	-0.02
30	Benzo(a)anthracene	0.400	0.400	0.0	93	-0.02
31	Chrysene	0.400	0.418	-4.5	97	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.260	35.0#	63	-0.02
33	Indeno(1,2,3-cd)pyrene	0.400	0.422	-5.5	96	-0.06
34 I	Perylene-d12	0.400	0.400	0.0	91	-0.03
35	Benzo(b)fluoranthene	0.400	0.424	-6.0	100	-0.03
36	Benzo(k)fluoranthene	0.400	0.437	-9.2	102	-0.03
37 C	Benzo(a)pyrene	0.400	0.408	-2.0	96	-0.03
38	Dibenzo(a,h)anthracene	0.400	0.427	-6.7	99	-0.07
39	Benzo(g,h,i)perylene	0.400	0.417	-4.2	97	-0.07

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

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