

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE101718\
 Data File : BE097966.D
 Acq On : 18 Oct 2018 2:53
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 18 10:19:56 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 15:36:44 2018
 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	114	0.00
2	1,4-Dioxane	0.400	0.402	-0.5	125	0.00
3	n-Nitrosodimethylamine	0.400	0.413	-3.2	134	0.00
4 S	2-Fluorophenol	0.400	0.420	-5.0	127	-0.01
5 S	Phenol-d6	0.400	0.443	-10.7	131	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.414	-3.5	127	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	127	-0.01
8 S	Nitrobenzene-d5	0.400	0.408	-2.0	133	-0.01
9	Naphthalene	0.400	0.436	-9.0	142	-0.01
10	Hexachlorobutadiene	0.400	0.431	-7.7	142	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.449	-12.2	147	-0.01
12	2-Methylnaphthalene	0.400	0.448	-12.0	148	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	134	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.406	-1.5	154	-0.01
15 S	2-Fluorobiphenyl	0.400	0.400	0.0	142	-0.01
16	Acenaphthylene	0.400	0.428	-7.0	151	0.00
17	Acenaphthene	0.400	0.424	-6.0	150	0.00
18	Fluorene	0.400	0.412	-3.0	145	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	128	-0.01
20	4-Bromophenyl-phenylether	0.400	0.436	-9.0	143	0.00
21	Hexachlorobenzene	0.400	0.454	-13.5	149	0.00
22	Pentachlorophenol	0.400	0.623	-55.7#	240	-0.01
23	Phenanthrene	0.400	0.423	-5.7	140	0.00
24	Anthracene	0.400	0.429	-7.2	144	-0.01
25 SURR	Fluoranthene-d10	0.400	0.405	-1.3	139	0.00
26	Fluoranthene	0.400	0.406	-1.5	139	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	119	-0.01
28	Pvrene	0.400	0.435	-8.7	138	0.00
29 S	Terphenyl-d14	0.400	0.418	-4.5	138	-0.01
30	Benzo(a)anthracene	0.400	0.424	-6.0	134	0.00
31	Chrysene	0.400	0.435	-8.7	142	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.455	-13.7	134	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.422	-5.5	128	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	118	-0.01
35	Benzo(b)fluoranthene	0.400	0.421	-5.2	133	-0.01
36	Benzo(k)fluoranthene	0.400	0.418	-4.5	136	0.00
37 C	Benzo(a)pyrene	0.400	0.426	-6.5	133	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.417	-4.2	131	-0.02
39	Benzo(a,h,i)perylene	0.400	0.404	-1.0	130	-0.02

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

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