

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE101818\
 Data File : BE097968.D
 Acq On : 18 Oct 2018 9:18
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 19 01:15:19 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	106	0.00
2	1,4-Dioxane	0.400	0.432	-8.0	124	0.00
3	n-Nitrosodimethylamine	0.400	0.394	1.5	118	0.00
4 S	2-Fluorophenol	0.400	0.436	-9.0	122	-0.01
5 S	Phenol-d6	0.400	0.479	-19.7	130	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.433	-8.2	122	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	121	-0.01
8 S	Nitrobenzene-d5	0.400	0.446	-11.5	137	-0.01
9	Naphthalene	0.400	0.449	-12.2	139	-0.01
10	Hexachlorobutadiene	0.400	0.449	-12.2	140	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.442	-10.5	137	-0.01
12	2-Methylnaphthalene	0.400	0.457	-14.2	143	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	127	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.403	-0.8	145	0.00
15 S	2-Fluorobiphenyl	0.400	0.419	-4.7	141	-0.01
16	Acenaphthylene	0.400	0.438	-9.5	147	0.00
17	Acenaphthene	0.400	0.435	-8.7	146	0.00
18	Fluorene	0.400	0.425	-6.2	142	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	124	-0.01
20	4-Bromophenyl-phenylether	0.400	0.439	-9.7	139	0.00
21	Hexachlorobenzene	0.400	0.465	-16.3	147	0.00
22	Pentachlorophenol	0.400	0.584	-46.0#	214	-0.01
23	Phenanthrene	0.400	0.425	-6.2	136	0.00
24	Anthracene	0.400	0.424	-6.0	137	0.00
25 SURR	Fluoranthene-d10	0.400	0.389	2.8	129	0.00
26	Fluoranthene	0.400	0.399	0.3	133	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	113	-0.01
28	Pyrene	0.400	0.443	-10.7	134	0.00
29 S	Terphenyl-d14	0.400	0.434	-8.5	136	-0.01
30	Benzo(a)anthracene	0.400	0.436	-9.0	131	0.00
31	Chrysene	0.400	0.445	-11.2	138	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.431	-7.7	122	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.440	-10.0	126	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	112	-0.01
35	Benzo(b)fluoranthene	0.400	0.429	-7.2	129	-0.01
36	Benzo(k)fluoranthene	0.400	0.430	-7.5	133	-0.01
37 C	Benzo(a)pyrene	0.400	0.433	-8.2	129	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.425	-6.2	127	-0.01
39	Benzo(g,h,i)perylene	0.400	0.416	-4.0	128	-0.02

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

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