

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE020419\
 Data File : BE098706.D
 Acq On : 4 Feb 2019 13:44
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 04 15:24:41 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE012419.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 24 14:44:44 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	108	0.00
2	1,4-Dioxane	0.400	0.337	15.8	102	0.00
3	n-Nitrosodimethylamine	0.400	0.349	12.8	97	0.00
4 S	2-Fluorophenol	0.400	0.338	15.5	91	0.00
5 S	Phenol-d6	0.400	0.357	10.8	102	0.00
6	bis(2-Chloroethyl)ether	0.400	0.306	23.5	91	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	109	0.00
8 S	Nitrobenzene-d5	0.400	0.386	3.5	105	0.00
9	Naphthalene	0.400	0.396	1.0	111	0.00
10	Hexachlorobutadiene	0.400	0.383	4.3	106	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.392	2.0	108	0.00
12	2-Methylnaphthalene	0.400	0.381	4.8	105	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	113	0.00
14 S	2,4,6-Tribromophenol	0.400	0.422	-5.5	116	-0.01
15 S	2-Fluorobiphenyl	0.400	0.386	3.5	110	0.00
16	Acenaphthylene	0.400	0.388	3.0	111	0.00
17	Acenaphthene	0.400	0.386	3.5	106	-0.02
18	Fluorene	0.400	0.389	2.8	111	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	111	0.00
20	4-Bromophenyl-phenylether	0.400	0.389	2.8	112	-0.01
21	Hexachlorobenzene	0.400	0.394	1.5	111	0.00
22	Pentachlorophenol	0.400	0.352	12.0	108	0.00
23	Phenanthrene	0.400	0.388	3.0	111	0.00
24	Anthracene	0.400	0.374	6.5	108	0.00
25 SURR	Fluoranthene-d10	0.400	0.383	4.3	109	0.00
26	Fluoranthene	0.400	0.388	3.0	111	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	108	-0.01
28	Pyrene	0.400	0.397	0.8	110	0.00
29 S	Terphenyl-d14	0.400	0.386	3.5	106	0.00
30	Benzo(a)anthracene	0.400	0.368	8.0	106	0.00
31	Chrysene	0.400	0.406	-1.5	113	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.373	6.8	106	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.394	1.5	116	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	113	0.00
35	Benzo(b)fluoranthene	0.400	0.381	4.8	109	0.00
36	Benzo(k)fluoranthene	0.400	0.371	7.3	108	0.00
37 C	Benzo(a)pyrene	0.400	0.393	1.8	115	0.00
38	Dibenzo(a,h)anthracene	0.400	0.374	6.5	115	0.00
39	Benzo(g,h,i)perylene	0.400	0.374	6.5	110	-0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 0