

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE111618\
 Data File : BE098242.D
 Acq On : 16 Nov 2018 22:50
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 16 23:20:43 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 16 19:05:33 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	108	0.00
2	1,4-Dioxane	0.400	0.372	7.0	106	0.00
3	n-Nitrosodimethylamine	0.400	0.419	-4.7	121	0.00
4 S	2-Fluorophenol	0.400	0.428	-7.0	115	0.00
5 S	Phenol-d6	0.400	0.450	-12.5	119	0.00
6	bis(2-Chloroethyl)ether	0.400	0.400	0.0	109	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	107	0.00
8 S	Nitrobenzene-d5	0.400	0.411	-2.7	108	0.00
9	Naphthalene	0.400	0.408	-2.0	106	0.00
10	Hexachlorobutadiene	0.400	0.405	-1.3	106	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	105	0.00
12	2-Methylnaphthalene	0.400	0.407	-1.7	108	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.400	0.520	-30.0#	139	-0.01
15 S	2-Fluorobiphenyl	0.400	0.387	3.3	102	0.00
16	Acenaphthylene	0.400	0.417	-4.2	106	0.00
17	Acenaphthene	0.400	0.428	-7.0	106	0.00
18	Fluorene	0.400	0.422	-5.5	108	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	104	0.00
20	4-Bromophenyl-phenylether	0.400	0.413	-3.2	107	0.00
21	Hexachlorobenzene	0.400	0.413	-3.2	110	-0.01
22	Pentachlorophenol	0.400	0.565	-41.2#	178	0.00
23	Phenanthrene	0.400	0.411	-2.7	107	-0.01
24	Anthracene	0.400	0.415	-3.7	109	0.00
25 SURR	Fluoranthene-d10	0.400	0.416	-4.0	109	0.00
26	Fluoranthene	0.400	0.407	-1.7	106	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	111	0.00
28	Pvrene	0.400	0.402	-0.5	108	0.00
29 S	Terphenyl-d14	0.400	0.407	-1.7	111	0.00
30	Benzo(a)anthracene	0.400	0.411	-2.7	114	0.00
31	Chrysene	0.400	0.411	-2.7	111	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.575	-43.7#	166	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.454	-13.5	123	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	113	0.00
35	Benzo(b)fluoranthene	0.400	0.415	-3.7	119	0.00
36	Benzo(k)fluoranthene	0.400	0.393	1.8	109	0.00
37 C	Benzo(a)pyrene	0.400	0.411	-2.7	116	0.00
38	Dibenzo(a,h)anthracene	0.400	0.465	-16.3	134	-0.01
39	Benzo(a,h,i)perylene	0.400	0.424	-6.0	126	0.00

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

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