

Data Path : Z:\HPCHEM1\BNA E\Data\BE032318\
 Data File : BE095830.D
 Acq On : 23 Mar 2018 19:25
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 23 23:23:58 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE032318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 23 12:41:11 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	107	0.00
2	1,4-Dioxane	0.400	0.367	8.3	103	0.00
3	n-Nitrosodimethylamine	0.400	0.338	15.5	87	0.00
4 S	2-Fluorophenol	0.400	0.354	11.5	91	0.00
5 S	Phenol-d6	0.400	0.368	8.0	94	0.00
6	bis(2-Chloroethyl)ether	0.400	0.370	7.5	95	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	106	0.00
8 S	Nitrobenzene-d5	0.400	0.364	9.0	96	0.00
9	Naphthalene	0.400	0.363	9.3	93	0.00
10	Hexachlorobutadiene	0.400	0.331	17.3	84	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.367	8.3	94	0.00
12	2-Methylnaphthalene	0.400	0.373	6.8	96	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	110	0.00
14 S	2,4,6-Tribromophenol	0.400	0.346	13.5	98	0.00
15 S	2-Fluorobiphenyl	0.400	0.356	11.0	95	-0.01
16	Acenaphthylene	0.400	0.349	12.8	95	0.00
17	Acenaphthene	0.400	0.350	12.5	96	0.00
18	Fluorene	0.400	0.350	12.5	95	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	111	0.00
20	4-Bromophenyl-phenylether	0.400	0.358	10.5	98	0.00
21	Hexachlorobenzene	0.400	0.355	11.3	97	0.00
22	Pentachlorophenol	0.400	0.320	20.0	82	0.00
23	Phenanthrene	0.400	0.364	9.0	98	0.00
24	Anthracene	0.400	0.363	9.3	100	0.00
25 SURR	Fluoranthene-d10	0.400	0.366	8.5	100	0.00
26	Fluoranthene	0.400	0.368	8.0	101	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	111	-0.01
28	Pvrene	0.400	0.380	5.0	96	-0.01
29 S	Terphenyl-d14	0.400	0.360	10.0	99	0.00
30	Benzo(a)anthracene	0.400	0.361	9.8	98	0.00
31	Chrysene	0.400	0.374	6.5	103	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.323	19.3	96	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.367	8.3	101	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	113	0.00
35	Benzo(b)fluoranthene	0.400	0.361	9.8	101	0.00
36	Benzo(k)fluoranthene	0.400	0.365	8.8	102	0.00
37 C	Benzo(a)pyrene	0.400	0.365	8.8	101	0.00
38	Dibenzo(a,h)anthracene	0.400	0.360	10.0	99	-0.01
39	Benzo(a,h,i)perylene	0.400	0.358	10.5	99	-0.01

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

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