

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051517\
 Data File : BE093033.D
 Acq On : 15 May 2017 14:23
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 16 01:10:38 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 15 13:35:57 2017
 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	100	0.00
2	1,4-Dioxane	0.400	0.350	12.5	101	0.00
3	n-Nitrosodimethylamine	0.400	0.358	10.5	102	0.00
4 S	2-Fluorophenol	0.400	0.352	12.0	99	0.00
5 S	Phenol-d6	0.400	0.358	10.5	102	0.00
6	bis(2-Chloroethyl)ether	0.400	0.363	9.3	103	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	103	0.00
8 S	Nitrobenzene-d5	0.400	0.343	14.2	102	0.00
9	Naphthalene	0.400	0.355	11.3	102	0.00
10	Hexachlorobutadiene	0.400	0.357	10.8	104	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.356	11.0	103	0.00
12	2-Methylnaphthalene	0.400	0.353	11.8	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	104	0.00
14 S	2,4,6-Tribromophenol	0.400	0.322	19.5	97	0.00
15 S	2-Fluorobiphenyl	0.400	0.354	11.5	102	0.00
16	Acenaphthylene	0.400	0.333	16.8	102	0.00
17	Acenaphthene	0.400	0.352	12.0	104	0.00
18	Fluorene	0.400	0.347	13.3	102	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	94	0.00
20	4-Bromophenyl-phenylether	0.400	0.335	16.3	98	0.00
21	Hexachlorobenzene	0.400	0.331	17.3	111	0.00
22	Pentachlorophenol	0.400	0.444	-11.0	148	0.00
23	Phenanthrene	0.400	0.325	18.8	99	0.03
24	Anthracene	0.400	0.320	20.0	109	0.00
25 SURR	Fluoranthene-d10	0.400	0.337	15.8	97	0.00
26	Fluoranthene	0.400	0.334	16.5	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	104	0.00
28	Pvrene	0.400	0.358	10.5	101	0.00
29 S	Terphenyl-d14	0.400	0.357	10.8	101	0.00
30	Benzo(a)anthracene	0.400	0.357	10.8	103	0.00
31	Chrysene	0.400	0.352	12.0	103	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.327	18.3	98	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.344	14.0	101	0.00
34 I	Perylene-d12	0.400	0.400	0.0	99	-0.01
35	Benzo(b)fluoranthene	0.400	0.343	14.2	99	0.00
36	Benzo(k)fluoranthene	0.400	0.347	13.3	97	0.00
37 C	Benzo(a)pyrene	0.400	0.352	12.0	102	0.00
38	Dibenzo(a,h)anthracene	0.400	0.352	12.0	100	0.00
39	Benzo(a,h,i)perylene	0.400	0.346	13.5	98	0.00

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

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