

Data Path : U:\HPCHEM1\BNA E\Data\BE082217\
 Data File : BE093871.D
 Acq On : 22 Aug 2017 12:18
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 23 01:20:14 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 14:30:21 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	78	0.00
2	1,4-Dioxane	0.400	0.308	23.0	59	0.00
3	n-Nitrosodimethylamine	0.400	0.346	13.5	66	0.01
4 S	2-Fluorophenol	0.400	0.344	14.0	68	0.00
5 S	Phenol-d6	0.400	0.339	15.3	68	0.01
6	bis(2-Chloroethyl)ether	0.400	0.343	14.2	69	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	79	0.00
8 S	Nitrobenzene-d5	0.400	0.354	11.5	73	0.00
9	Naphthalene	0.400	0.352	12.0	70	0.00
10	Hexachlorobutadiene	0.400	0.330	17.5	66	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.351	12.3	69	0.00
12	2-Methylnaphthalene	0.400	0.346	13.5	69	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	77	0.00
14 S	2,4,6-Tribromophenol	0.400	0.321	19.8	68	0.03
15 S	2-Fluorobiphenyl	0.400	0.350	12.5	67	0.00
16	Acenaphthylene	0.400	0.348	13.0	69	0.00
17	Acenaphthene	0.400	0.340	15.0	66	0.00
18	Fluorene	0.400	0.344	14.0	67	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	71	0.00
20	4-Bromophenyl-phenylether	0.400	0.460	-15.0	83	0.00
21	Hexachlorobenzene	0.400	0.410	-2.5	70	0.00
22	Pentachlorophenol	0.400	0.260	35.0#	49	0.00
23	Phenanthrene	0.400	0.380	5.0	68	0.00
24	Anthracene	0.400	0.344	14.0	62	0.00
25 SURR	Fluoranthene-d10	0.400	0.374	6.5	62	0.00
26	Fluoranthene	0.400	0.363	9.3	61	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	68	0.01
28	Pvrene	0.400	0.349	12.8	61	0.00
29 S	Terphenyl-d14	0.400	0.348	13.0	61	0.00
30	Benzo(a)anthracene	0.400	0.332	17.0	58	0.00
31	Chrysene	0.400	0.357	10.8	62	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.357	10.8	60	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.286	28.5#	49	0.04
34 I	Perylene-d12	0.400	0.400	0.0	67	0.01
35	Benzo(b)fluoranthene	0.400	0.310	22.5	51	0.01
36	Benzo(k)fluoranthene	0.400	0.358	10.5	62	0.01
37 C	Benzo(a)pyrene	0.400	0.335	16.3	56	0.02
38	Dibenzo(a,h)anthracene	0.400	0.292	27.0#	50	0.04
39	Benzo(a,h,i)perylene	0.400	0.297	25.8#	51	0.04

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

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