

Data Path : Z:\HPCHEM1\BNA_E\Data\BE090117\
 Data File : BE093901.D
 Acq On : 1 Sep 2017 14:28
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 02 01:18:42 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE082617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 01 16:31:36 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	118	0.00
2	1,4-Dioxane	0.400	0.347	13.3	110	0.00
3	n-Nitrosodimethylamine	0.400	0.387	3.3	121	-0.01
4 S	2-Fluorophenol	0.400	0.380	5.0	115	0.00
5 S	Phenol-d6	0.400	0.367	8.3	114	0.00
6	bis(2-Chloroethyl)ether	0.400	0.375	6.3	115	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	124	0.00
8 S	Nitrobenzene-d5	0.400	0.371	7.3	120	0.00
9	Naphthalene	0.400	0.345	13.8	108	0.00
10	Hexachlorobutadiene	0.400	0.368	8.0	118	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.351	12.3	111	0.00
12	2-Methylnaphthalene	0.400	0.342	14.5	109	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	124	0.00
14 S	2,4,6-Tribromophenol	0.400	0.346	13.5	125	0.00
15 S	2-Fluorobiphenyl	0.400	0.342	14.5	110	-0.01
16	Acenaphthylene	0.400	0.354	11.5	118	0.00
17	Acenaphthene	0.400	0.346	13.5	113	0.00
18	Fluorene	0.400	0.354	11.5	116	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	129	0.00
20	4-Bromophenyl-phenylether	0.400	0.346	13.5	109	0.00
21	Hexachlorobenzene	0.400	0.363	9.3	109	0.00
22	Pentachlorophenol	0.400	0.367	8.3	146	0.00
23	Phenanthrene	0.400	0.320	20.0	107	0.00
24	Anthracene	0.400	0.347	13.3	121	0.00
25 SURR	Fluoranthene-d10	0.400	0.404	-1.0	125	0.00
26	Fluoranthene	0.400	0.400	0.0	125	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	133	0.00
28	Pyrene	0.400	0.358	10.5	127	0.00
29 S	Terphenyl-d14	0.400	0.358	10.5	126	0.00
30	Benzo(a)anthracene	0.400	0.367	8.3	132	-0.01
31	Chrysene	0.400	0.343	14.2	120	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.457	-14.2	171	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.398	0.5	142	0.00
34 I	Perylene-d12	0.400	0.400	0.0	144	0.00
35	Benzo(b)fluoranthene	0.400	0.349	12.8	138	0.00
36	Benzo(k)fluoranthene	0.400	0.323	19.3	120	0.00
37 C	Benzo(a)pyrene	0.400	0.347	13.3	134	0.00
38	Dibenzo(a,h)anthracene	0.400	0.375	6.3	142	0.00
39	Benzo(g,h,i)perylene	0.400	0.380	5.0	147	0.00

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

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