

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051617\
 Data File : BE093047.D
 Acq On : 16 May 2017 10:28
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 18 00:53:08 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	99	0.00
2	1,4-Dioxane	0.400	0.356	11.0	102	0.00
3	n-Nitrosodimethylamine	0.400	0.366	8.5	104	-0.01
4 S	2-Fluorophenol	0.400	0.343	14.2	96	0.00
5 S	Phenol-d6	0.400	0.364	9.0	103	0.00
6	bis(2-Chloroethyl)ether	0.400	0.367	8.3	104	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	0.00
8 S	Nitrobenzene-d5	0.400	0.347	13.3	101	0.00
9	Naphthalene	0.400	0.353	11.8	99	0.00
10	Hexachlorobutadiene	0.400	0.361	9.8	103	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.358	10.5	100	0.01
12	2-Methylnaphthalene	0.400	0.355	11.3	101	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.400	0.367	8.3	104	0.00
15 S	2-Fluorobiphenyl	0.400	0.376	6.0	102	0.00
16	Acenaphthylene	0.400	0.350	12.5	100	0.00
17	Acenaphthene	0.400	0.358	10.5	99	0.02
18	Fluorene	0.400	0.364	9.0	100	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	99	0.00
20	4-Bromophenyl-phenylether	0.400	0.324	19.0	101	0.00
21	Hexachlorobenzene	0.400	0.296	26.0#	108	0.00
22	Pentachlorophenol	0.400	0.369	7.8	131	0.00
23	Phenanthrene	0.400	0.299	25.3#	98	0.03
24	Anthracene	0.400	0.294	26.5#	108	0.00
25 SURR	Fluoranthene-d10	0.400	0.308	23.0	94	0.00
26	Fluoranthene	0.400	0.315	21.3	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pvrene	0.400	0.382	4.5	102	0.00
29 S	Terphenyl-d14	0.400	0.377	5.8	102	0.00
30	Benzo(a)anthracene	0.400	0.358	10.5	99	0.00
31	Chrysene	0.400	0.360	10.0	100	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.353	11.8	102	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.377	5.8	105	0.00
34 I	Perylene-d12	0.400	0.400	0.0	101	-0.01
35	Benzo(b)fluoranthene	0.400	0.348	13.0	103	0.00
36	Benzo(k)fluoranthene	0.400	0.343	14.2	98	0.00
37 C	Benzo(a)pyrene	0.400	0.347	13.3	103	0.00
38	Dibenzo(a,h)anthracene	0.400	0.361	9.8	105	0.00
39	Benzo(a,h,i)perylene	0.400	0.350	12.5	101	0.00

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051617\
Data File : BE093047.D
Acq On : 16 May 2017 10:28
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: May 18 00:53:08 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)

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