

Data Path : Z:\HPCHEM1\BNA E\Data\BE100417\
 Data File : BE094170.D
 Acq On : 4 Oct 2017 17:08
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 04 20:21:15 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 29 13:57:50 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	110	0.00
2	1,4-Dioxane	0.400	0.372	7.0	99	0.00
3	n-Nitrosodimethylamine	0.400	0.296	26.0#	82	0.00
4 S	2-Fluorophenol	0.400	0.353	11.8	96	0.00
5 S	Phenol-d6	0.400	0.399	0.3	109	0.00
6	bis(2-Chloroethyl)ether	0.400	0.371	7.3	100	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	127	0.00
8 S	Nitrobenzene-d5	0.400	0.319	20.3	107	0.00
9	Naphthalene	0.400	0.402	-0.5	129	0.00
10	Hexachlorobutadiene	0.400	0.375	6.3	125	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.334	16.5	108	0.00
12	2-Methylnaphthalene	0.400	0.336	16.0	109	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	115	0.00
14 S	2,4,6-Tribromophenol	0.400	0.252	37.0#	80	0.00
15 S	2-Fluorobiphenyl	0.400	0.390	2.5	108	0.00
16	Acenaphthylene	0.400	0.395	1.3	111	0.00
17	Acenaphthene	0.400	0.423	-5.7	114	0.00
18	Fluorene	0.400	0.415	-3.7	116	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.400	0.525	-31.3#	121	0.00
21	Hexachlorobenzene	0.400	0.723	-80.7#	164	0.00
22	Pentachlorophenol	0.400	0.229	42.8#	53	0.00
23	Phenanthrene	0.400	0.520	-30.0#	122	0.00
24	Anthracene	0.400	0.392	2.0	93	0.00
25 SURR	Fluoranthene-d10	0.400	0.394	1.5	93	0.00
26	Fluoranthene	0.400	0.408	-2.0	94	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	81	0.00
28	Pvrene	0.400	0.432	-8.0	91	0.00
29 S	Terphenyl-d14	0.400	0.410	-2.5	87	0.00
30	Benzo(a)anthracene	0.400	0.386	3.5	78	0.00
31	Chrysene	0.400	0.392	2.0	79	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.089	77.8#	49	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.375	6.3	79	0.00
34 I	Perylene-d12	0.400	0.400	0.0	79	0.00
35	Benzo(b)fluoranthene	0.400	0.355	11.3	78	0.00
36	Benzo(k)fluoranthene	0.400	0.390	2.5	77	0.00
37 C	Benzo(a)pyrene	0.400	0.391	2.3	80	0.00
38	Dibenzo(a,h)anthracene	0.400	0.375	6.3	78	0.00
39	Benzo(a,h,i)perylene	0.400	0.387	3.3	79	0.00

Data Path : Z:\HPCHEM1\BNA E\Data\BE100417\
Data File : BE094170.D
Acq On : 4 Oct 2017 17:08
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 04 20:21:15 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 29 13:57:50 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				