

Data Path : Z:\HPCHEM1\BNA E\Data\BE092517\
 Data File : BE094108.D
 Acq On : 25 Sep 2017 18:39
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 26 02:22:55 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 25 17:37:08 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	99	0.00
2	1,4-Dioxane	0.780	0.792	-1.5	97	0.00
3	n-Nitrosodimethylamine	0.996	0.844	15.3	84	0.00
4 S	2-Fluorophenol	1.516	1.440	5.0	93	0.00
5 S	Phenol-d6	2.106	2.025	3.8	94	0.00
6	bis(2-Chloroethyl)ether	1.632	1.592	2.5	94	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
8 S	Nitrobenzene-d5	0.387	0.354	8.5	92	0.00
9	Naphthalene	4.516	4.501	0.3	95	0.00
10	Hexachlorobutadiene	0.169	0.167	1.2	98	0.00
11 SURR	2-Methylnaphthalene-d10	0.616	0.607	1.5	95	0.00
12	2-Methylnaphthalene	0.731	0.729	0.3	97	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.256	0.191	25.4#	86	0.00
15 S	2-Fluorobiphenyl	1.322	1.352	-2.3	96	0.01
16	Acenaphthylene	2.005	1.965	2.0	94	0.00
17	Acenaphthene	1.253	1.305	-4.2	96	0.00
18	Fluorene	1.700	1.737	-2.2	98	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.164	0.160	2.4	95	0.00
21	Hexachlorobenzene	0.104	0.111	-6.7	103	0.00
22	Pentachlorophenol	0.095	0.079	16.8	82	0.00
23	Phenanthrene	1.023	1.062	-3.8	102	0.00
24	Anthracene	1.074	1.041	3.1	97	0.00
25 SURR	Fluoranthene-d10	3.873	3.903	-0.8	100	0.00
26	Fluoranthene	1.147	1.181	-3.0	101	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	102	0.00
28	Pvrene	1.267	1.187	6.3	100	0.00
29 S	Terphenyl-d14	0.730	0.671	8.1	99	0.00
30	Benzo(a)anthracene	1.347	1.291	4.2	98	0.00
31	Chrysene	1.300	1.281	1.5	101	0.00
32	Bis(2-ethylhexyl)phthalate	6.447	4.686	27.3#	100	0.00
33	Indeno(1,2,3-cd)pyrene	1.391	1.295	6.9	99	0.00
34 I	Perylene-d12	1.000	1.000	0.0	100	0.00
35	Benzo(b)fluoranthene	1.443	1.315	8.9	101	0.00
36	Benzo(k)fluoranthene	1.382	1.299	6.0	94	0.00
37 C	Benzo(a)pyrene	1.304	1.246	4.4	99	0.00
38	Dibenzo(a,h)anthracene	1.235	1.170	5.3	99	0.00
39	Benzo(a,h,i)perylene	1.222	1.163	4.8	99	0.00

Data Path : Z:\HPCHEM1\BNA E\Data\BE092517\
Data File : BE094108.D
Acq On : 25 Sep 2017 18:39
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Sep 26 02:22:55 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Sep 25 17:37:08 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	AvgRF	CCRF	%Dev Area	% Dev(min)

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