

Data Path : Z:\HPCHEM1\BNA_E\Data\BE101617\
 Data File : BE094424.D
 Acq On : 16 Oct 2017 19:52
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 16 20:34:53 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 14:54:40 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	93	-0.02
2	1,4-Dioxane	0.698	0.681	2.4	90	0.00
3	n-Nitrosodimethylamine	0.695	0.702	-1.0	94	0.00
4 S	2-Fluorophenol	1.352	1.368	-1.2	92	0.00
5 S	Phenol-d6	2.108	2.073	1.7	88	0.00
6	bis(2-Chloroethyl)ether	1.559	1.575	-1.0	90	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
8 S	Nitrobenzene-d5	0.348	0.336	3.4	89	-0.02
9	Naphthalene	4.473	4.586	-2.5	94	-0.02
10	Hexachlorobutadiene	0.179	0.175	2.2	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.617	0.637	-3.2	95	0.00
12	2-Methylnaphthalene	0.737	0.783	-6.2	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
14 S	2,4,6-Tribromophenol	0.152	0.153	-0.7	94	0.00
15 S	2-Fluorobiphenyl	1.428	1.422	0.4	97	0.00
16	Acenaphthylene	2.169	2.159	0.5	99	0.00
17	Acenaphthene	1.366	1.369	-0.2	101	0.00
18	Fluorene	1.761	1.730	1.8	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
20	4-Bromophenyl-phenylether	0.200	0.221	-10.5	105	-0.03
21	Hexachlorobenzene	0.365	0.406	-11.2	98	0.00
22	Pentachlorophenol	0.032	0.031	3.1	89	0.00
23	Phenanthrene	1.574	1.492	5.2	78	0.00
24	Anthracene	1.216	1.310	-7.7	96	0.00
25 SURR	Fluoranthene-d10	6.441	6.956	-8.0	96	0.00
26	Fluoranthene	1.954	2.109	-7.9	97	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
28	Pyrene	1.509	1.435	4.9	97	0.00
29 S	Terphenyl-d14	0.807	0.782	3.1	97	0.00
30	Benzo(a)anthracene	1.384	1.355	2.1	100	-0.01
31	Chrysene	1.332	1.306	2.0	97	-0.01
32	Bis(2-ethylhexyl)phthalate	3.984	3.701	7.1	98	0.00
33	Indeno(1,2,3-cd)pyrene	1.341	1.304	2.8	99	0.00
34 I	Perylene-d12	1.000	1.000	0.0	99	0.00
35	Benzo(b)fluoranthene	1.357	1.353	0.3	99	0.00
36	Benzo(k)fluoranthene	1.337	1.384	-3.5	101	0.00
37 C	Benzo(a)pyrene	1.275	1.280	-0.4	99	0.00
38	Dibenzo(a,h)anthracene	1.207	1.216	-0.7	99	0.00
39	Benzo(g,h,i)perylene	1.161	1.191	-2.6	101	-0.01

Data Path : Z:\HPCHEM1\BNA_E\Data\BE101617\
Data File : BE094424.D
Acq On : 16 Oct 2017 19:52
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 16 20:34:53 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE101617.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 16 14:54:40 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			