

Data Path : Z:\HPCHEM1\BNA_E\Data\BE121517\
 Data File : BE094995.D
 Acq On : 15 Dec 2017 11:03
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 15 16:27:35 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 14 18:42:53 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	98	0.00
2	1,4-Dioxane	0.394	0.384	2.5	90	0.00
3	n-Nitrosodimethylamine	0.466	0.474	-1.7	99	0.00
4 S	2-Fluorophenol	0.621	0.655	-5.5	107	0.00
5 S	Phenol-d6	0.923	0.984	-6.6	107	0.00
6	bis(2-Chloroethyl)ether	0.901	0.880	2.3	96	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8 S	Nitrobenzene-d5	0.328	0.338	-3.0	108	0.00
9	Naphthalene	4.697	4.622	1.6	99	0.00
10	Hexachlorobutadiene	0.214	0.205	4.2	96	0.00
11 SURR	2-Methylnaphthalene-d10	0.655	0.646	1.4	100	0.00
12	2-Methylnaphthalene	1.172	1.136	3.1	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
14 S	2,4,6-Tribromophenol	0.093	0.110	-18.3	132	0.00
15 S	2-Fluorobiphenyl	1.772	1.656	6.5	96	0.00
16	Acenaphthylene	2.138	2.036	4.8	104	0.00
17	Acenaphthene	1.418	1.364	3.8	101	0.00
18	Fluorene	1.762	1.724	2.2	104	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.234	0.226	3.4	102	0.00
21	Hexachlorobenzene	0.239	0.238	0.4	103	0.00
22	Pentachlorophenol	0.046	0.065	-41.3#	150#	0.00
23	Phenanthrene	1.246	1.220	2.1	103	0.00
24	Anthracene	1.279	1.248	2.4	106	0.00
25 SURR	Fluoranthene-d10	5.189	4.962	4.4	102	0.00
26	Fluoranthene	1.538	1.480	3.8	102	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	100	0.00
28	Pyrene	1.355	1.565	-15.5	112	0.00
29 S	Terphenyl-d14	0.753	0.770	-2.3	102	0.00
30	Benzo(a)anthracene	1.169	1.173	-0.3	102	0.00
31	Chrysene	1.316	1.255	4.6	96	0.00
32	Bis(2-ethylhexyl)phthalate	1.489	1.740	-16.9	131	-0.01
33	Indeno(1,2,3-cd)pyrene	1.071	1.112	-3.8	102	0.00
34 I	Perylene-d12	1.000	1.000	0.0	102	0.00
35	Benzo(b)fluoranthene	1.469	1.431	2.6	99	0.00
36	Benzo(k)fluoranthene	1.488	1.348	9.4	97	0.00
37 C	Benzo(a)pyrene	1.303	1.255	3.7	99	0.00
38	Dibenzo(a,h)anthracene	1.209	1.176	2.7	100	0.00
39	Benzo(g,h,i)perylene	1.203	1.152	4.2	100	0.00

Data Path : Z:\HPCHEM1\BNA_E\Data\BE121517\
Data File : BE094995.D
Acq On : 15 Dec 2017 11:03
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Dec 15 16:27:35 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 14 18:42:53 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		