

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE112118\
 Data File : BE098254.D
 Acq On : 21 Nov 2018 9:38
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 21 10:48:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 16 19:05:33 2018
 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	108	0.00
2	1,4-Dioxane	0.628	0.643	-2.4	116	0.00
3	n-Nitrosodimethylamine	0.706	0.708	-0.3	116	0.00
4 S	2-Fluorophenol	1.073	1.042	2.9	104	0.00
5 S	Phenol-d6	1.685	1.755	-4.2	109	0.01
6	bis(2-Chloroethyl)ether	1.443	1.379	4.4	104	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
8 S	Nitrobenzene-d5	0.378	0.392	-3.7	109	0.00
9	Naphthalene	4.073	4.072	0.0	104	0.00
10	Hexachlorobutadiene	0.239	0.222	7.1	97	0.00
11 SURR	2-Methylnaphthalene-d10	0.702	0.673	4.1	101	0.00
12	2-Methylnaphthalene	0.718	0.700	2.5	103	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
14 S	2,4,6-Tribromophenol	0.129	0.117	9.3	102	0.00
15 S	2-Fluorobiphenyl	1.739	1.630	6.3	104	0.00
16	Acenaphthylene	1.904	1.866	2.0	105	0.00
17	Acenaphthene	1.137	1.130	0.6	104	0.00
18	Fluorene	1.449	1.444	0.3	107	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
20	4-Bromophenyl-phenylether	0.248	0.238	4.0	104	0.00
21	Hexachlorobenzene	0.256	0.242	5.5	105	-0.01
22	Pentachlorophenol	0.035	0.041	-17.1	131	0.00
23	Phenanthrene	1.029	1.006	2.2	106	-0.01
24	Anthracene	0.935	0.908	2.9	106	0.00
25 SURR	Fluoranthene-d10	4.866	4.690	3.6	106	0.00
26	Fluoranthene	1.235	1.195	3.2	105	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	112	0.00
28	Pvrene	1.169	1.129	3.4	104	0.00
29 S	Terphenyl-d14	0.919	0.870	5.3	103	0.00
30	Benzo(a)anthracene	1.157	1.095	5.4	105	0.00
31	Chrysene	1.191	1.187	0.3	109	0.00
32	Bis(2-ethylhexyl)phthalate	1.804	1.719	4.7	110	0.00
33	Indeno(1,2,3-cd)pyrene	1.019	1.084	-6.4	116	0.00
34 I	Perylene-d12	1.000	1.000	0.0	113	0.00
35	Benzo(b)fluoranthene	1.130	1.089	3.6	110	0.00
36	Benzo(k)fluoranthene	1.276	1.198	6.1	103	0.00
37 C	Benzo(a)pyrene	1.109	1.080	2.6	109	0.00
38	Dibenzo(a,h)anthracene	0.889	0.893	-0.4	115	0.00
39	Benzo(a,h,i)perylene	0.958	0.959	-0.1	118	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE112118\
Data File : BE098254.D
Acq On : 21 Nov 2018 9:38
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Nov 21 10:48:38 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE111618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 16 19:05:33 2018
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			