

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE012020\
 Data File : BE101102.D
 Acq On : 20 Jan 2020 15:05
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 21 10:41:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 18:36:25 2020
 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	52	-0.01
2	1,4-Dioxane	1.020	1.036	-1.6	57	0.00
3	n-Nitrosodimethylamine	0.556	0.502	9.7	54	-0.01
4 S	2-Fluorophenol	0.900	1.036	-15.1	60	-0.01
5 S	Phenol-d6	1.139	1.069	6.1	54	-0.01
6	bis(2-Chloroethyl)ether	1.357	1.113	18.0	46#	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	54	0.00
8 S	Nitrobenzene-d5	0.286	0.259	9.4	54	0.00
9	Naphthalene	4.441	4.394	1.1	52	0.00
10	Hexachlorobutadiene	0.188	0.199	-5.9	55	-0.01
11 SURR	2-Methylnaphthalene-d10	0.763	0.754	1.2	53	0.00
12	2-Methylnaphthalene	0.672	0.633	5.8	52	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	64	0.00
14 S	2,4,6-Tribromophenol	0.113	0.097	14.2	63	0.00
15 S	2-Fluorobiphenyl	1.519	1.304	14.2	54	-0.01
16	Acenaphthylene	1.629	1.537	5.6	67	0.00
17	Acenaphthene	1.166	1.072	8.1	61	-0.01
18	Fluorene	1.474	1.393	5.5	62	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	72	0.00
20	4-Bromophenyl-phenylether	0.195	0.178	8.7	69	-0.01
21	Hexachlorobenzene	0.214	0.193	9.8	64	-0.01
22	Pentachlorophenol	0.053	0.024	54.7#	43#	0.00
23	Phenanthrene	1.092	0.953	12.7	64	0.00
24	Anthracene	1.031	1.015	1.6	75	0.00
25 SURR	Fluoranthene-d10	5.555	5.841	-5.1	78	-0.01
26	Fluoranthene	1.368	1.417	-3.6	80	-0.01
27 I	Chrysene-d12	1.000	1.000	0.0	99	-0.01
28	Pvrene	1.489	1.254	15.8	81	0.00
29 S	Terphenyl-d14	0.982	0.820	16.5	81	-0.01
30	Benzo(a)anthracene	1.141	1.106	3.1	100	0.00
31	Chrysene	1.696	1.681	0.9	95	-0.01
32	Bis(2-ethylhexyl)phthalate	1.902	2.658	-39.7#	158#	0.00
33	Indeno(1,2,3-cd)pyrene	1.340	1.448	-8.1	109	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	115	-0.01
35	Benzo(b)fluoranthene	1.122	0.934	16.8	103	-0.01
36	Benzo(k)fluoranthene	1.626	1.433	11.9	102	-0.01
37 C	Benzo(a)pyrene	1.085	1.025	5.5	113	-0.01
38	Dibenzo(a,h)anthracene	0.998	0.922	7.6	116	-0.01
39	Benzo(a,h,i)perylene	1.227	1.070	12.8	98	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE012020\
Data File : BE101102.D
Acq On : 20 Jan 2020 15:05
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jan 21 10:41:02 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 08 18:36:25 2020
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			