

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102319\
 Data File : BE100662.D
 Acq On : 23 Oct 2019 15:27
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 23 16:58:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 16:49:33 2019
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	54	0.00
2	1,4-Dioxane	0.400	0.404	-1.0	55	0.00
3	n-Nitrosodimethylamine	0.400	0.393	1.8	58	0.00
4 S	2-Fluorophenol	0.400	0.429	-7.2	62	0.00
5 S	Phenol-d6	0.400	0.438	-9.5	63	0.00
6	bis(2-Chloroethyl)ether	0.400	0.388	3.0	56	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	57	0.00
8 S	Nitrobenzene-d5	0.400	0.563	-40.7#	91	0.00
9	Naphthalene	0.400	0.389	2.8	57	0.00
10	Hexachlorobutadiene	0.400	0.400	0.0	59	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	58	0.00
12	2-Methylnaphthalene	0.400	0.398	0.5	59	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	59	0.00
14 S	2,4,6-Tribromophenol	0.400	0.647	-61.7#	109	0.00
15 S	2-Fluorobiphenyl	0.400	0.427	-6.7	65	0.00
16	Acenaphthylene	0.400	0.402	-0.5	62	0.00
17	Acenaphthene	0.400	0.390	2.5	60	0.00
18	Fluorene	0.400	0.394	1.5	60	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	64	0.00
20	4-Bromophenyl-phenylether	0.400	0.395	1.3	65	0.00
21	Hexachlorobenzene	0.400	0.395	1.3	67	0.00
22	Pentachlorophenol	0.400	0.648	-62.0#	117	0.00
23	Phenanthrene	0.400	0.408	-2.0	65	0.00
24	Anthracene	0.400	0.381	4.8	63	0.00
25 SURR	Fluoranthene-d10	0.400	0.414	-3.5	69	0.00
26	Fluoranthene	0.400	0.430	-7.5	69	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	71	0.00
28	Pvrene	0.400	0.400	0.0	71	0.00
29 S	Terphenyl-d14	0.400	0.434	-8.5	81	0.00
30	Benzo(a)anthracene	0.400	0.387	3.3	69	0.00
31	Chrysene	0.400	0.393	1.8	70	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.454	-13.5	80	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.370	7.5	68	0.00
34 I	Perylene-d12	0.400	0.400	0.0	64	0.00
35	Benzo(b)fluoranthene	0.400	0.385	3.8	67	0.00
36	Benzo(k)fluoranthene	0.400	0.382	4.5	66	0.00
37 C	Benzo(a)pyrene	0.400	0.380	5.0	66	0.00
38	Dibenzo(a,h)anthracene	0.400	0.371	7.3	66	0.00
39	Benzo(a,h,i)perylene	0.400	0.383	4.3	66	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102319\
Data File : BE100662.D
Acq On : 23 Oct 2019 15:27
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Oct 23 16:58:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 23 16:49:33 2019
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	Amount	Calc.	%Dev	Area	% Dev(min)

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