

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011320\
 Data File : BE101045.D
 Acq On : 13 Jan 2020 19:26
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 14 09:58:34 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	92	0.00
2	1,4-Dioxane	0.400	0.429	-7.2	99	0.00
3	n-Nitrosodimethylamine	0.400	0.267	33.3#	71	0.00
4 S	2-Fluorophenol	0.400	0.346	13.5	81	0.00
5 S	Phenol-d6	0.400	0.335	16.3	87	0.00
6	bis(2-Chloroethyl)ether	0.400	0.311	22.3	78	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	104	0.00
8 S	Nitrobenzene-d5	0.400	0.277	30.8#	81	0.00
9	Naphthalene	0.400	0.396	1.0	101	0.00
10	Hexachlorobutadiene	0.400	0.408	-2.0	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	103	0.00
12	2-Methylnaphthalene	0.400	0.374	6.5	99	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.400	0.266	33.5#	79	0.00
15 S	2-Fluorobiphenyl	0.400	0.412	-3.0	104	-0.01
16	Acenaphthylene	0.400	0.328	18.0	93	0.00
17	Acenaphthene	0.400	0.401	-0.3	106	-0.01
18	Fluorene	0.400	0.391	2.3	102	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20	4-Bromophenyl-phenylether	0.400	0.365	8.8	99	0.00
21	Hexachlorobenzene	0.400	0.404	-1.0	103	-0.01
22	Pentachlorophenol	0.400	0.411	-2.7	77	0.00
23	Phenanthrene	0.400	0.387	3.3	101	0.00
24	Anthracene	0.400	0.361	9.8	98	0.00
25 SURR	Fluoranthene-d10	0.400	0.365	8.8	97	0.00
26	Fluoranthene	0.400	0.360	10.0	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	97	-0.01
28	Pvrene	0.400	0.407	-1.7	96	0.00
29 S	Terphenyl-d14	0.400	0.401	-0.3	96	0.00
30	Benzo(a)anthracene	0.400	0.346	13.5	87	0.00
31	Chrysene	0.400	0.432	-8.0	102	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.299	25.3#	83	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.418	-4.5	104	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	108	0.00
35	Benzo(b)fluoranthene	0.400	0.336	16.0	97	0.00
36	Benzo(k)fluoranthene	0.400	0.355	11.3	96	0.00
37 C	Benzo(a)pyrene	0.400	0.380	5.0	106	0.00
38	Dibenzo(a,h)anthracene	0.400	0.375	6.3	110	-0.01
39	Benzo(a,h,i)perylene	0.400	0.380	5.0	100	-0.02

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011320\
Data File : BE101045.D
Acq On : 13 Jan 2020 19:26
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jan 14 09:58:34 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	Amount	Calc.	%Dev	Area	% Dev(min)

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