

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011020\
 Data File : BE101024.D
 Acq On : 10 Jan 2020 15:02
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 10 15:59:13 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	103	0.00
2	1,4-Dioxane	0.400	0.340	15.0	94	0.00
3	n-Nitrosodimethylamine	0.400	0.351	12.3	105	0.00
4 S	2-Fluorophenol	0.400	0.348	13.0	90	0.00
5 S	Phenol-d6	0.400	0.328	18.0	95	0.00
6	bis(2-Chloroethyl)ether	0.400	0.330	17.5	93	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	106	0.00
8 S	Nitrobenzene-d5	0.400	0.300	25.0#	89	0.00
9	Naphthalene	0.400	0.407	-1.7	106	0.00
10	Hexachlorobutadiene	0.400	0.406	-1.5	104	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	108	0.00
12	2-Methylnaphthalene	0.400	0.388	3.0	105	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	104	0.00
14 S	2,4,6-Tribromophenol	0.400	0.286	28.5#	86	0.00
15 S	2-Fluorobiphenyl	0.400	0.407	-1.7	105	0.00
16	Acenaphthylene	0.400	0.346	13.5	100	0.00
17	Acenaphthene	0.400	0.396	1.0	106	0.00
18	Fluorene	0.400	0.390	2.5	104	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.400	0.378	5.5	101	0.00
21	Hexachlorobenzene	0.400	0.424	-6.0	107	0.00
22	Pentachlorophenol	0.400	0.401	-0.3	72	0.00
23	Phenanthrene	0.400	0.395	1.3	102	0.00
24	Anthracene	0.400	0.377	5.8	101	0.00
25 SURR	Fluoranthene-d10	0.400	0.360	10.0	95	0.00
26	Fluoranthene	0.400	0.363	9.3	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	89	0.00
28	Pvrene	0.400	0.434	-8.5	94	0.00
29 S	Terphenyl-d14	0.400	0.425	-6.2	93	0.00
30	Benzo(a)anthracene	0.400	0.311	22.3	72	0.00
31	Chrysene	0.400	0.458	-14.5	98	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.293	26.8#	74	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.401	-0.3	91	0.00
34 I	Perylene-d12	0.400	0.400	0.0	94	0.00
35	Benzo(b)fluoranthene	0.400	0.320	20.0	80	0.00
36	Benzo(k)fluoranthene	0.400	0.430	-7.5	101	0.00
37 C	Benzo(a)pyrene	0.400	0.375	6.3	91	0.00
38	Dibenzo(a,h)anthracene	0.400	0.354	11.5	90	-0.01
39	Benzo(a,h,i)perylene	0.400	0.378	5.5	86	-0.01

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011020\
Data File : BE101024.D
Acq On : 10 Jan 2020 15:02
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_E
LabSampleId :
SSTDCCC0.4

Quant Time: Jan 10 15:59:13 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				