

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011420\
 Data File : BE101047.D
 Acq On : 14 Jan 2020 9:37
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 15 10:03:58 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	98	-0.01
2	1,4-Dioxane	0.400	0.422	-5.5	104	0.00
3	n-Nitrosodimethylamine	0.400	0.295	26.3#	84	0.01
4 S	2-Fluorophenol	0.400	0.348	13.0	86	0.00
5 S	Phenol-d6	0.400	0.362	9.5	100	0.00
6	bis(2-Chloroethyl)ether	0.400	0.405	-1.3	109	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	106	0.00
8 S	Nitrobenzene-d5	0.400	0.324	19.0	96	0.01
9	Naphthalene	0.400	0.410	-2.5	106	0.00
10	Hexachlorobutadiene	0.400	0.429	-7.2	109	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.412	-3.0	109	0.00
12	2-Methylnaphthalene	0.400	0.396	1.0	106	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	108	0.00
14 S	2,4,6-Tribromophenol	0.400	0.225	43.8#	70	0.00
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	109	0.00
16	Acenaphthylene	0.400	0.315	21.3	94	0.00
17	Acenaphthene	0.400	0.394	1.5	109	-0.01
18	Fluorene	0.400	0.384	4.0	106	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	104	0.00
20	4-Bromophenyl-phenylether	0.400	0.369	7.8	101	0.00
21	Hexachlorobenzene	0.400	0.424	-6.0	109	0.00
22	Pentachlorophenol	0.400	0.383	4.3	64	0.00
23	Phenanthrene	0.400	0.393	1.8	103	0.00
24	Anthracene	0.400	0.364	9.0	100	0.00
25 SURR	Fluoranthene-d10	0.400	0.358	10.5	96	0.00
26	Fluoranthene	0.400	0.359	10.3	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	89	-0.01
28	Pvrene	0.400	0.437	-9.2	94	0.00
29 S	Terphenyl-d14	0.400	0.428	-7.0	94	0.00
30	Benzo(a)anthracene	0.400	0.307	23.3	71	0.00
31	Chrysene	0.400	0.449	-12.2	97	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.273	31.8#	70	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.409	-2.2	93	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	98	0.00
35	Benzo(b)fluoranthene	0.400	0.328	18.0	86	0.00
36	Benzo(k)fluoranthene	0.400	0.381	4.8	93	0.00
37 C	Benzo(a)pyrene	0.400	0.364	9.0	92	0.00
38	Dibenzo(a,h)anthracene	0.400	0.354	11.5	94	-0.01
39	Benzo(a,h,i)perylene	0.400	0.381	4.8	91	-0.01

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Compound	Amount	Calc.	%Dev	Area	% Dev(min)

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