

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011120\
 Data File : BE101033.D
 Acq On : 10 Jan 2020 21:20
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 11 01:48:55 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	97	0.00
2	1,4-Dioxane	0.400	0.412	-3.0	101	0.00
3	n-Nitrosodimethylamine	0.400	0.310	22.5	87	0.00
4 S	2-Fluorophenol	0.400	0.352	12.0	86	0.00
5 S	Phenol-d6	0.400	0.327	18.3	89	0.00
6	bis(2-Chloroethyl)ether	0.400	0.306	23.5	81	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	108	0.00
8 S	Nitrobenzene-d5	0.400	0.228	43.0#	69	0.01
9	Naphthalene	0.400	0.391	2.3	103	0.00
10	Hexachlorobutadiene	0.400	0.398	0.5	103	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.392	2.0	105	0.00
12	2-Methylnaphthalene	0.400	0.376	6.0	103	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	105	0.00
14 S	2,4,6-Tribromophenol	0.400	0.239	40.3#	73	0.00
15 S	2-Fluorobiphenyl	0.400	0.398	0.5	104	0.00
16	Acenaphthylene	0.400	0.311	22.3	91	0.00
17	Acenaphthene	0.400	0.392	2.0	106	0.00
18	Fluorene	0.400	0.386	3.5	104	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	102	0.00
20	4-Bromophenyl-phenylether	0.400	0.365	8.8	98	0.00
21	Hexachlorobenzene	0.400	0.396	1.0	100	0.00
22	Pentachlorophenol	0.400	0.386	3.5	64	0.00
23	Phenanthrene	0.400	0.378	5.5	98	0.00
24	Anthracene	0.400	0.362	9.5	97	0.00
25 SURR	Fluoranthene-d10	0.400	0.359	10.3	95	0.00
26	Fluoranthene	0.400	0.356	11.0	97	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	114	-0.01
28	Pvrene	0.400	0.338	15.5	94	0.00
29 S	Terphenyl-d14	0.400	0.332	17.0	93	0.00
30	Benzo(a)anthracene	0.400	0.345	13.8	102	0.00
31	Chrysene	0.400	0.445	-11.2	123	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.497	-24.2	162	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.462	-15.5	135	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	127	-0.01
35	Benzo(b)fluoranthene	0.400	0.324	19.0	110	-0.01
36	Benzo(k)fluoranthene	0.400	0.369	7.8	117	0.00
37 C	Benzo(a)pyrene	0.400	0.398	0.5	130	0.00
38	Dibenzo(a,h)anthracene	0.400	0.356	11.0	122	-0.02
39	Benzo(a,h,i)perylene	0.400	0.383	4.3	118	-0.02

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

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