

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE012120\
 Data File : BE101104.D
 Acq On : 21 Jan 2020 9:36
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 22 07:55:16 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	70	-0.01
2	1,4-Dioxane	0.400	0.347	13.3	64	0.00
3	n-Nitrosodimethylamine	0.400	0.335	16.3	68	-0.01
4 S	2-Fluorophenol	0.400	0.411	-2.7	72	-0.01
5 S	Phenol-d6	0.400	0.348	13.0	68	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.311	22.3	59	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	71	0.00
8 S	Nitrobenzene-d5	0.400	0.348	13.0	69	0.00
9	Naphthalene	0.400	0.391	2.3	68	0.00
10	Hexachlorobutadiene	0.400	0.428	-7.0	73	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	71	0.00
12	2-Methylnaphthalene	0.400	0.368	8.0	66	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	83	0.00
14 S	2,4,6-Tribromophenol	0.400	0.293	26.8#	70	0.00
15 S	2-Fluorobiphenyl	0.400	0.351	12.3	72	0.00
16	Acenaphthylene	0.400	0.361	9.8	83	0.00
17	Acenaphthene	0.400	0.367	8.3	78	-0.01
18	Fluorene	0.400	0.385	3.8	81	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.400	0.379	5.3	91	0.00
21	Hexachlorobenzene	0.400	0.358	10.5	81	-0.01
22	Pentachlorophenol	0.400	0.339	15.3	38	0.00
23	Phenanthrene	0.400	0.363	9.3	84	0.00
24	Anthracene	0.400	0.380	5.0	92	0.00
25 SURR	Fluoranthene-d10	0.400	0.432	-8.0	103	0.00
26	Fluoranthene	0.400	0.425	-6.2	104	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	115	-0.01
28	Pvrene	0.400	0.372	7.0	104	0.00
29 S	Terphenyl-d14	0.400	0.378	5.5	107	0.00
30	Benzo(a)anthracene	0.400	0.344	14.0	103	0.00
31	Chrysene	0.400	0.401	-0.3	112	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.290	27.5#	96	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.342	14.5	101	0.00
34 I	Perylene-d12	0.400	0.400	0.0	103	-0.01
35	Benzo(b)fluoranthene	0.400	0.330	17.5	91	0.00
36	Benzo(k)fluoranthene	0.400	0.335	16.3	86	0.00
37 C	Benzo(a)pyrene	0.400	0.377	5.8	100	0.00
38	Dibenzo(a,h)anthracene	0.400	0.380	5.0	106	-0.02
39	Benzo(a,h,i)perylene	0.400	0.367	8.3	92	-0.01

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

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