

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011720\
 Data File : BE101086.D
 Acq On : 17 Jan 2020 15:53
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 17 23:25:50 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	48	0.00
2	1,4-Dioxane	0.400	0.451	-12.7	54	0.00
3	n-Nitrosodimethylamine	0.400	0.425	-6.2	59	0.00
4 S	2-Fluorophenol	0.400	0.505	-26.2#	61	0.00
5 S	Phenol-d6	0.400	0.398	0.5	54	0.00
6	bis(2-Chloroethyl)ether	0.400	0.384	4.0	51	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	52	0.00
8 S	Nitrobenzene-d5	0.400	0.364	9.0	53	0.00
9	Naphthalene	0.400	0.402	-0.5	51	0.00
10	Hexachlorobutadiene	0.400	0.416	-4.0	52	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	52	-0.01
12	2-Methylnaphthalene	0.400	0.401	-0.3	53	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	67	0.00
14 S	2,4,6-Tribromophenol	0.400	0.371	7.3	71	0.00
15 S	2-Fluorobiphenyl	0.400	0.344	14.0	57	-0.01
16	Acenaphthylene	0.400	0.356	11.0	66	-0.01
17	Acenaphthene	0.400	0.374	6.5	64	-0.01
18	Fluorene	0.400	0.402	-0.5	68	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.400	0.313	21.8	72	-0.01
21	Hexachlorobenzene	0.400	0.342	14.5	74	-0.01
22	Pentachlorophenol	0.400	0.384	4.0	54	-0.01
23	Phenanthrene	0.400	0.361	9.8	80	0.00
24	Anthracene	0.400	0.388	3.0	89	0.00
25 SURR	Fluoranthene-d10	0.400	0.464	-16.0	105	-0.01
26	Fluoranthene	0.400	0.478	-19.5	111	-0.01
27 I	Chrysene-d12	0.400	0.400	0.0	146	-0.01
28	Pvrene	0.400	0.323	19.3	115	0.00
29 S	Terphenyl-d14	0.400	0.361	9.8	130	-0.01
30	Benzo(a)anthracene	0.400	0.409	-2.2	155	0.00
31	Chrysene	0.400	0.404	-1.0	143	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.466	-16.5	195	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.288	28.0#	108	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	119	-0.01
35	Benzo(b)fluoranthene	0.400	0.382	4.5	122	-0.01
36	Benzo(k)fluoranthene	0.400	0.358	10.5	107	-0.01
37 C	Benzo(a)pyrene	0.400	0.398	0.5	122	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.331	17.3	107	-0.02
39	Benzo(a,h,i)perylene	0.400	0.319	20.3	93	-0.01

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

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