

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN042120\
 Data File : BN010560.D
 Acq On : 21 Apr 2020 11:07
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 21 13:14:29 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN041320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 13 18:04:22 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	76	0.00
2	1,4-Dioxane	0.400	0.389	2.8	75	0.00
3	n-Nitrosodimethylamine	0.400	0.406	-1.5	92	0.00
4 S	2-Fluorophenol	0.400	0.422	-5.5	80	0.00
5 S	Phenol-d6	0.400	0.402	-0.5	76	0.00
6	bis(2-Chloroethyl)ether	0.400	0.417	-4.2	78	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	78	0.00
8 S	Nitrobenzene-d5	0.400	0.388	3.0	80	0.00
9	Naphthalene	0.400	0.407	-1.7	81	0.00
10	Hexachlorobutadiene	0.400	0.410	-2.5	80	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.459	-14.8	91	0.00
12	2-Methylnaphthalene	0.400	0.410	-2.5	81	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	80	0.00
14 S	2,4,6-Tribromophenol	0.400	0.317	20.8	67	0.00
15 S	2-Fluorobiphenyl	0.400	0.405	-1.3	81	0.00
16	Acenaphthylene	0.400	0.371	7.3	76	0.00
17	Acenaphthene	0.400	0.402	-0.5	82	0.00
18	Fluorene	0.400	0.400	0.0	81	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	80	0.00
20	4-Bromophenyl-phenylether	0.400	0.390	2.5	81	0.00
21	Hexachlorobenzene	0.400	0.426	-6.5	85	0.00
22	Pentachlorophenol	0.400	0.287	28.3#	68	0.00
23	Phenanthrene	0.400	0.407	-1.7	82	0.00
24	Anthracene	0.400	0.368	8.0	76	0.00
25 SURR	Fluoranthene-d10	0.400	0.431	-7.7	88	0.00
26	Fluoranthene	0.400	0.393	1.8	79	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	72	0.00
28	Pvrene	0.400	0.424	-6.0	77	0.00
29 S	Terphenyl-d14	0.400	0.425	-6.2	78	0.00
30	Benzo(a)anthracene	0.400	0.361	9.8	65	0.00
31	Chrysene	0.400	0.405	-1.3	74	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.272	32.0#	52	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.432	-8.0	77	0.00
34 I	Perylene-d12	0.400	0.400	0.0	71	0.00
35	Benzo(b)fluoranthene	0.400	0.407	-1.7	73	0.00
36	Benzo(k)fluoranthene	0.400	0.410	-2.5	74	0.00
37 C	Benzo(a)pyrene	0.400	0.390	2.5	70	0.00
38	Dibenzo(a,h)anthracene	0.400	0.424	-6.0	76	0.00
39	Benzo(a,h,i)perylene	0.400	0.444	-11.0	79	0.00

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

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