

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN022120\
 Data File : BN009804.D
 Acq On : 21 Feb 2020 12:36
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 21 16:37:16 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN021920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 19 15:57:43 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	123	-0.02
2	1,4-Dioxane	0.400	0.358	10.5	111	-0.02
3	n-Nitrosodimethylamine	0.400	0.376	6.0	132	-0.02
4 S	2-Fluorophenol	0.400	0.414	-3.5	132	-0.02
5 S	Phenol-d6	0.400	0.358	10.5	116	-0.02
6	bis(2-Chloroethyl)ether	0.400	0.403	-0.8	129	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	128	-0.03
8 S	Nitrobenzene-d5	0.400	0.347	13.3	118	-0.01
9	Naphthalene	0.400	0.392	2.0	129	-0.03
10	Hexachlorobutadiene	0.400	0.397	0.8	127	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.386	3.5	128	-0.02
12	2-Methylnaphthalene	0.400	0.385	3.8	128	-0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	131	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.391	2.3	143	-0.01
15 S	2-Fluorobiphenyl	0.400	0.372	7.0	127	-0.02
16	Acenaphthylene	0.400	0.380	5.0	139	-0.02
17	Acenaphthene	0.400	0.387	3.3	132	-0.02
18	Fluorene	0.400	0.373	6.8	126	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	123	-0.01
20	4-Bromophenyl-phenylether	0.400	0.214	46.5#	84	-0.01
21	Hexachlorobenzene	0.400	0.404	-1.0	130	-0.02
22	Pentachlorophenol	0.400	0.432	-8.0	137	-0.02
23	Phenanthrene	0.400	0.372	7.0	124	-0.01
24	Anthracene	0.400	0.371	7.3	127	-0.01
25 SURR	Fluoranthene-d10	0.400	0.370	7.5	122	-0.02
26	Fluoranthene	0.400	0.369	7.8	122	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	120	-0.02
28	Pvrene	0.400	0.403	-0.8	125	-0.02
29 S	Terphenyl-d14	0.400	0.381	4.8	118	-0.02
30	Benzo(a)anthracene	0.400	0.372	7.0	126	-0.02
31	Chrysene	0.400	0.363	9.3	112	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.436	-9.0	146	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	118	-0.03
34 I	Perylene-d12	0.400	0.400	0.0	125	-0.02
35	Benzo(b)fluoranthene	0.400	0.338	15.5	115	-0.02
36	Benzo(k)fluoranthene	0.400	0.386	3.5	127	-0.02
37 C	Benzo(a)pyrene	0.400	0.376	6.0	121	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.347	13.3	115	-0.03
39	Benzo(a,h,i)perylene	0.400	0.371	7.3	123	-0.03

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

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