

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031020\
 Data File : BN010174.D
 Acq On : 11 Mar 2020 03:55
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 11 04:31:09 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	84	0.00
2	1,4-Dioxane	0.400	0.412	-3.0	87	0.00
3	n-Nitrosodimethylamine	0.400	0.529	-32.3#	126	0.02
4 S	2-Fluorophenol	0.400	0.398	0.5	86	0.00
5 S	Phenol-d6	0.400	0.403	-0.8	89	0.02
6	bis(2-Chloroethyl)ether	0.400	0.498	-24.5	108	0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	76	0.01
8 S	Nitrobenzene-d5	0.400	0.379	5.3	76	0.03
9	Naphthalene	0.400	0.414	-3.5	81	0.01
10	Hexachlorobutadiene	0.400	0.449	-12.2	85	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.431	-7.7	85	0.01
12	2-Methylnaphthalene	0.400	0.445	-11.2	88	0.03
13 I	Acenaphthene-d10	0.400	0.400	0.0	74	0.00
14 S	2,4,6-Tribromophenol	0.400	0.387	3.3	80	0.01
15 S	2-Fluorobiphenyl	0.400	0.442	-10.5	85	0.01
16	Acenaphthylene	0.400	0.391	2.3	81	0.00
17	Acenaphthene	0.400	0.399	0.3	77	0.00
18	Fluorene	0.400	0.379	5.3	72	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	66	0.01
20	4-Bromophenyl-phenylether	0.400	0.253	36.8#	50	0.00
21	Hexachlorobenzene	0.400	0.427	-6.7	73	0.01
22	Pentachlorophenol	0.400	0.454	-13.5	79	0.01
23	Phenanthrene	0.400	0.386	3.5	68	0.01
24	Anthracene	0.400	0.386	3.5	70	0.01
25 SURR	Fluoranthene-d10	0.400	0.397	0.8	70	0.00
26	Fluoranthene	0.400	0.398	0.5	70	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	72	0.00
28	Pvrene	0.400	0.432	-8.0	80	0.00
29 S	Terphenyl-d14	0.400	0.406	-1.5	75	0.01
30	Benzo(a)anthracene	0.400	0.354	11.5	72	0.00
31	Chrysene	0.400	0.422	-5.5	78	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.449	-12.2	90	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.452	-13.0	81	0.00
34 I	Perylene-d12	0.400	0.400	0.0	74	0.00
35	Benzo(b)fluoranthene	0.400	0.388	3.0	79	0.00
36	Benzo(k)fluoranthene	0.400	0.407	-1.7	80	0.00
37 C	Benzo(a)pyrene	0.400	0.398	0.5	76	0.00
38	Dibenzo(a,h)anthracene	0.400	0.406	-1.5	80	0.01
39	Benzo(a,h,i)perylene	0.400	0.426	-6.5	84	0.00

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

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