

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN073018\
 Data File : BN002190.D
 Acq On : 30 Jul 2018 17:50
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 31 01:11:09 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 16:17:29 2018
 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	87	0.00
2	1,4-Dioxane	0.400	0.386	3.5	93	0.00
3	n-Nitrosodimethylamine	0.400	0.355	11.3	80	0.00
4 S	2-Fluorophenol	0.400	0.375	6.3	86	0.00
5 S	Phenol-d6	0.400	0.373	6.8	87	0.00
6	bis(2-Chloroethyl)ether	0.400	0.372	7.0	81	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	90	0.00
8 S	Nitrobenzene-d5	0.400	0.382	4.5	88	0.00
9	Naphthalene	0.400	0.391	2.3	90	0.00
10	Hexachlorobutadiene	0.400	0.389	2.8	90	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.385	3.8	90	0.00
12	2-Methylnaphthalene	0.400	0.388	3.0	89	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	85	0.00
14 S	2,4,6-Tribromophenol	0.400	0.335	16.3	72	0.00
15 S	2-Fluorobiphenyl	0.400	0.394	1.5	88	0.00
16	Acenaphthylene	0.400	0.376	6.0	83	0.01
17	Acenaphthene	0.400	0.388	3.0	85	0.00
18	Fluorene	0.400	0.380	5.0	81	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	71	0.00
20	4-Bromophenyl-phenylether	0.400	0.395	1.3	76	0.00
21	Hexachlorobenzene	0.400	0.408	-2.0	77	0.00
22	Pentachlorophenol	0.400	0.342	14.5	63	0.00
23	Phenanthrene	0.400	0.382	4.5	70	0.00
24	Anthracene	0.400	0.363	9.3	68	0.00
25 SURR	Fluoranthene-d10	0.400	0.347	13.3	60	0.00
26	Fluoranthene	0.400	0.345	13.8	59	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	54	0.00
28	Pvrene	0.400	0.410	-2.5	59	0.00
29 S	Terphenyl-d14	0.400	0.395	1.3	57	0.00
30	Benzo(a)anthracene	0.400	0.369	7.8	52	0.00
31	Chrysene	0.400	0.389	2.8	54	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.336	16.0	46	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.435	-8.7	69	0.00
34 I	Perylene-d12	0.400	0.400	0.0	60	0.00
35	Benzo(b)fluoranthene	0.400	0.367	8.3	56	0.00
36	Benzo(k)fluoranthene	0.400	0.385	3.8	61	0.00
37 C	Benzo(a)pyrene	0.400	0.389	2.8	61	0.00
38	Dibenzo(a,h)anthracene	0.400	0.404	-1.0	69	0.00
39	Benzo(a,h,i)perylene	0.400	0.409	-2.2	69	0.00

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

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