

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN083019\
 Data File : BN007520.D
 Acq On : 30 Aug 2019 13:03
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 31 01:58:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN082319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 29 12:18:45 2019
 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	184	-0.02
2	1,4-Dioxane	0.400	0.438	-9.5	194	0.00
3	n-Nitrosodimethylamine	0.400	0.328	18.0	145	0.00
4 S	2-Fluorophenol	0.400	0.360	10.0	164	-0.02
5 S	Phenol-d6	0.400	0.356	11.0	171	0.00
6	bis(2-Chloroethyl)ether	0.400	0.420	-5.0	195	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	202	-0.03
8 S	Nitrobenzene-d5	0.400	0.381	4.8	201	-0.01
9	Naphthalene	0.400	0.393	1.8	193	-0.01
10	Hexachlorobutadiene	0.400	0.359	10.3	174	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.411	-2.7	204	-0.01
12	2-Methylnaphthalene	0.400	0.419	-4.7	205	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	234	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.459	-14.8	274	-0.02
15 S	2-Fluorobiphenyl	0.400	0.386	3.5	221	0.00
16	Acenaphthylene	0.400	0.372	7.0	215	-0.02
17	Acenaphthene	0.400	0.388	3.0	223	0.00
18	Fluorene	0.400	0.405	-1.3	232	-0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	244	0.00
20	4-Bromophenyl-phenylether	0.400	0.113	71.8#	79	0.00
21	Hexachlorobenzene	0.400	0.390	2.5	236	-0.02
22	Pentachlorophenol	0.400	0.604	-51.0#	394	-0.02
23	Phenanthrene	0.400	0.410	-2.5	250	-0.02
24	Anthracene	0.400	0.395	1.3	239	0.00
25 SURR	Fluoranthene-d10	0.400	0.386	3.5	228	-0.01
26	Fluoranthene	0.400	0.393	1.8	234	-0.01
27 I	Chrysene-d12	0.400	0.400	0.0	217	0.00
28	Pvrene	0.400	0.438	-9.5	232	0.00
29 S	Terphenyl-d14	0.400	0.433	-8.2	227	-0.01
30	Benzo(a)anthracene	0.400	0.402	-0.5	214	-0.01
31	Chrysene	0.400	0.395	1.3	205	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.507	-26.7#	261	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.314	21.5	165	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	185	-0.01
35	Benzo(b)fluoranthene	0.400	0.418	-4.5	189	0.00
36	Benzo(k)fluoranthene	0.400	0.431	-7.7	192	0.00
37 C	Benzo(a)pyrene	0.400	0.405	-1.3	182	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.354	11.5	159	-0.02
39	Benzo(a,h,i)perylene	0.400	0.363	9.3	163	-0.02

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

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