

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090819\
 Data File : BN007559.D
 Acq On : 08 Sep 2019 20:24
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 09 04:46:46 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	312	-0.02
2	1,4-Dioxane	0.400	0.396	1.0	297	0.00
3	n-Nitrosodimethylamine	0.400	0.303	24.3	228	0.00
4 S	2-Fluorophenol	0.400	0.332	17.0	258	-0.02
5 S	Phenol-d6	0.400	0.328	18.0	269	0.00
6	bis(2-Chloroethyl)ether	0.400	0.424	-6.0	333	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	310	-0.03
8 S	Nitrobenzene-d5	0.400	0.371	7.3	301	-0.01
9	Naphthalene	0.400	0.396	1.0	300	-0.01
10	Hexachlorobutadiene	0.400	0.347	13.3	259	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.368	8.0	281	-0.01
12	2-Methylnaphthalene	0.400	0.377	5.8	284	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	302	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.283	29.3#	218	-0.02
15 S	2-Fluorobiphenyl	0.400	0.391	2.3	289	0.00
16	Acenaphthylene	0.400	0.344	14.0	256	-0.02
17	Acenaphthene	0.400	0.390	2.5	288	0.00
18	Fluorene	0.400	0.379	5.3	280	-0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	287	-0.02
20	4-Bromophenyl-phenylether	0.400	0.142	64.5#	118	0.00
21	Hexachlorobenzene	0.400	0.385	3.8	273	-0.02
22	Pentachlorophenol	0.400	0.233	41.8#	178	-0.02
23	Phenanthrene	0.400	0.406	-1.5	291	-0.02
24	Anthracene	0.400	0.352	12.0	251	0.00
25 SURR	Fluoranthene-d10	0.400	0.352	12.0	244	-0.02
26	Fluoranthene	0.400	0.361	9.8	252	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	276	-0.01
28	Pvrene	0.400	0.370	7.5	249	-0.01
29 S	Terphenyl-d14	0.400	0.360	10.0	240	-0.02
30	Benzo(a)anthracene	0.400	0.342	14.5	232	-0.01
31	Chrysene	0.400	0.388	3.0	257	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.230	42.5#	157	-0.02
33	Indeno(1,2,3-cd)pyrene	0.400	0.361	9.8	240	-0.04
34 I	Perylene-d12	0.400	0.400	0.0	266	-0.02
35	Benzo(b)fluoranthene	0.400	0.374	6.5	243	-0.02
36	Benzo(k)fluoranthene	0.400	0.426	-6.5	272	-0.02
37 C	Benzo(a)pyrene	0.400	0.367	8.3	237	-0.03
38	Dibenzo(a,h)anthracene	0.400	0.364	9.0	234	-0.04
39	Benzo(a,h,i)perylene	0.400	0.365	8.8	236	-0.04

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

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