

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE010419\
 Data File : BE098619.D
 Acq On : 4 Jan 2019 17:27
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 04 21:14:23 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 02 10:30:27 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	70	0.00
2	1,4-Dioxane	0.400	0.408	-2.0	73	0.00
3	n-Nitrosodimethylamine	0.400	0.403	-0.8	77	0.01
4 S	2-Fluorophenol	0.400	0.539	-34.8#	98	0.01
5 S	Phenol-d6	0.400	0.565	-41.2#	99	0.01
6	bis(2-Chloroethyl)ether	0.400	0.422	-5.5	77	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	74	0.00
8 S	Nitrobenzene-d5	0.400	0.419	-4.7	84	0.00
9	Naphthalene	0.400	0.413	-3.2	78	0.00
10	Hexachlorobutadiene	0.400	0.463	-15.8	88	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.425	-6.2	79	0.00
12	2-Methylnaphthalene	0.400	0.428	-7.0	80	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	74	0.00
14 S	2,4,6-Tribromophenol	0.400	0.517	-29.3#	101	0.01
15 S	2-Fluorobiphenyl	0.400	0.392	2.0	80	0.00
16	Acenaphthylene	0.400	0.389	2.8	75	0.02
17	Acenaphthene	0.400	0.405	-1.3	76	0.02
18	Fluorene	0.400	0.396	1.0	74	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	62	0.00
20	4-Bromophenyl-phenylether	0.400	0.424	-6.0	69	0.01
21	Hexachlorobenzene	0.400	0.444	-11.0	71	0.01
22	Pentachlorophenol	0.400	0.815	-103.7#	149	0.01
23	Phenanthrene	0.400	0.408	-2.0	64	0.01
24	Anthracene	0.400	0.393	1.8	62	0.01
25 SURR	Fluoranthene-d10	0.400	0.402	-0.5	60	0.00
26	Fluoranthene	0.400	0.400	0.0	60	0.02
27 I	Chrysene-d12	0.400	0.400	0.0	68	0.01
28	Pvrene	0.400	0.372	7.0	61	0.02
29 S	Terphenyl-d14	0.400	0.382	4.5	65	0.00
30	Benzo(a)anthracene	0.400	0.416	-4.0	72	0.00
31	Chrysene	0.400	0.389	2.8	66	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.490	-22.5	83	0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.431	-7.7	77	-0.06
34 I	Perylene-d12	0.400	0.400	0.0	73	-0.03
35	Benzo(b)fluoranthene	0.400	0.418	-4.5	79	-0.01
36	Benzo(k)fluoranthene	0.400	0.400	0.0	76	-0.01
37 C	Benzo(a)pyrene	0.400	0.411	-2.7	77	-0.03
38	Dibenzo(a,h)anthracene	0.400	0.404	-1.0	76	-0.06
39	Benzo(a,h,i)perylene	0.400	0.410	-2.5	78	-0.08

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

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