

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE071719\
 Data File : BE100449.D
 Acq On : 17 Jul 2019 14:02
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 17 15:25:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 16 14:03:26 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	83	0.00
2	1,4-Dioxane	0.400	0.408	-2.0	89	0.00
3	n-Nitrosodimethylamine	0.400	0.394	1.5	94	0.00
4 S	2-Fluorophenol	0.400	0.491	-22.7	107	0.00
5 S	Phenol-d6	0.400	0.465	-16.3	101	0.00
6	bis(2-Chloroethyl)ether	0.400	0.413	-3.2	87	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	87	0.00
8 S	Nitrobenzene-d5	0.400	0.321	19.8	78	0.00
9	Naphthalene	0.400	0.391	2.3	87	0.00
10	Hexachlorobutadiene	0.400	0.384	4.0	86	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	94	0.00
12	2-Methylnaphthalene	0.400	0.408	-2.0	94	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	107	0.00
14 S	2,4,6-Tribromophenol	0.400	0.434	-8.5	132	0.00
15 S	2-Fluorobiphenyl	0.400	0.357	10.8	96	0.00
16	Acenaphthylene	0.400	0.432	-8.0	127	0.00
17	Acenaphthene	0.400	0.384	4.0	109	0.00
18	Fluorene	0.400	0.400	0.0	112	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	112	0.00
20	4-Bromophenyl-phenylether	0.400	0.355	11.3	106	0.01
21	Hexachlorobenzene	0.400	0.322	19.5	92	0.00
22	Pentachlorophenol	0.400	0.499	-24.7	138	0.00
23	Phenanthrene	0.400	0.361	9.8	104	0.00
24	Anthracene	0.400	0.402	-0.5	127	0.00
25 SURR	Fluoranthene-d10	0.400	0.427	-6.7	127	0.00
26	Fluoranthene	0.400	0.395	1.3	117	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	110	0.00
28	Pvrene	0.400	0.428	-7.0	125	0.00
29 S	Terphenyl-d14	0.400	0.393	1.8	114	0.00
30	Benzo(a)anthracene	0.400	0.448	-12.0	139	0.00
31	Chrysene	0.400	0.379	5.3	108	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	1.069	-167.2#	351	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.423	-5.7	130	0.00
34 I	Perylene-d12	0.400	0.400	0.0	128	0.00
35	Benzo(b)fluoranthene	0.400	0.358	10.5	123	0.00
36	Benzo(k)fluoranthene	0.400	0.357	10.8	124	0.00
37 C	Benzo(a)pyrene	0.400	0.424	-6.0	155	0.00
38	Dibenzo(a,h)anthracene	0.400	0.348	13.0	124	0.00
39	Benzo(a,h,i)perylene	0.400	0.345	13.8	118	0.00

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

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