

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE072319\
 Data File : BE100504.D
 Acq On : 23 Jul 2019 16:47
 Operator : HP/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 24 02:07:34 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE072219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 23 08:06:59 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	109	0.00
2	1,4-Dioxane	0.400	0.393	1.8	115	0.00
3	n-Nitrosodimethylamine	0.400	0.497	-24.2	136	-0.01
4 S	2-Fluorophenol	0.400	0.451	-12.7	126	0.00
5 S	Phenol-d6	0.400	0.439	-9.7	122	0.00
6	bis(2-Chloroethyl)ether	0.400	0.408	-2.0	113	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	103	-0.01
8 S	Nitrobenzene-d5	0.400	0.461	-15.3	126	0.00
9	Naphthalene	0.400	0.390	2.5	101	0.00
10	Hexachlorobutadiene	0.400	0.390	2.5	100	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.378	5.5	100	0.00
12	2-Methylnaphthalene	0.400	0.375	6.3	99	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	94	0.00
14 S	2,4,6-Tribromophenol	0.400	0.479	-19.7	131	0.00
15 S	2-Fluorobiphenyl	0.400	0.386	3.5	97	0.00
16	Acenaphthylene	0.400	0.392	2.0	100	0.00
17	Acenaphthene	0.400	0.379	5.3	95	0.00
18	Fluorene	0.400	0.389	2.8	96	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	96	0.00
20	4-Bromophenyl-phenylether	0.400	0.387	3.3	96	-0.01
21	Hexachlorobenzene	0.400	0.373	6.8	93	-0.01
22	Pentachlorophenol	0.400	0.512	-28.0#	128	-0.01
23	Phenanthrene	0.400	0.392	2.0	97	-0.01
24	Anthracene	0.400	0.390	2.5	98	0.00
25 SURR	Fluoranthene-d10	0.400	0.370	7.5	92	0.00
26	Fluoranthene	0.400	0.373	6.8	91	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	87	0.00
28	Pvrene	0.400	0.410	-2.5	91	0.00
29 S	Terphenyl-d14	0.400	0.402	-0.5	90	0.00
30	Benzo(a)anthracene	0.400	0.390	2.5	87	-0.01
31	Chrysene	0.400	0.406	-1.5	90	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.420	-5.0	92	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.412	-3.0	91	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	90	-0.01
35	Benzo(b)fluoranthene	0.400	0.391	2.3	92	0.00
36	Benzo(k)fluoranthene	0.400	0.385	3.8	89	0.00
37 C	Benzo(a)pyrene	0.400	0.391	2.3	92	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.391	2.3	93	-0.01
39	Benzo(a,h,i)perylene	0.400	0.388	3.0	91	-0.02

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

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