

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102319\
 Data File : BE100673.D
 Acq On : 23 Oct 2019 22:09
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 24 01:34:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 16:49:33 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	59	0.00
2	1,4-Dioxane	0.400	0.350	12.5	52	0.00
3	n-Nitrosodimethylamine	0.400	0.333	16.8	53	0.00
4 S	2-Fluorophenol	0.400	0.400	0.0	62	0.00
5 S	Phenol-d6	0.400	0.407	-1.7	63	0.00
6	bis(2-Chloroethyl)ether	0.400	0.362	9.5	57	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	60	0.00
8 S	Nitrobenzene-d5	0.400	0.528	-32.0#	91	0.00
9	Naphthalene	0.400	0.378	5.5	58	0.00
10	Hexachlorobutadiene	0.400	0.388	3.0	61	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.376	6.0	59	0.00
12	2-Methylnaphthalene	0.400	0.371	7.3	58	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	63	0.00
14 S	2,4,6-Tribromophenol	0.400	0.528	-32.0#	94	0.00
15 S	2-Fluorobiphenyl	0.400	0.413	-3.2	66	0.00
16	Acenaphthylene	0.400	0.360	10.0	59	-0.01
17	Acenaphthene	0.400	0.366	8.5	60	0.00
18	Fluorene	0.400	0.366	8.5	60	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	67	-0.01
20	4-Bromophenyl-phenylether	0.400	0.365	8.8	63	0.00
21	Hexachlorobenzene	0.400	0.375	6.3	66	-0.01
22	Pentachlorophenol	0.400	0.497	-24.2	94	0.00
23	Phenanthrene	0.400	0.386	3.5	64	-0.01
24	Anthracene	0.400	0.342	14.5	59	0.00
25 SURR	Fluoranthene-d10	0.400	0.376	6.0	65	0.00
26	Fluoranthene	0.400	0.396	1.0	66	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	70	0.00
28	Pvrene	0.400	0.380	5.0	67	0.00
29 S	Terphenyl-d14	0.400	0.413	-3.2	76	0.00
30	Benzo(a)anthracene	0.400	0.362	9.5	64	0.00
31	Chrysene	0.400	0.382	4.5	67	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.361	9.8	63	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.360	10.0	65	0.00
34 I	Perylene-d12	0.400	0.400	0.0	68	0.00
35	Benzo(b)fluoranthene	0.400	0.354	11.5	65	0.00
36	Benzo(k)fluoranthene	0.400	0.364	9.0	66	0.00
37 C	Benzo(a)pyrene	0.400	0.351	12.3	64	0.00
38	Dibenzo(a,h)anthracene	0.400	0.350	12.5	65	0.00
39	Benzo(a,h,i)perylene	0.400	0.358	10.5	65	0.00

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

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