

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE011320\
 Data File : BE101035.D
 Acq On : 13 Jan 2020 10:53
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 13 14:29:23 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE010820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 08 18:36:25 2020
 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	103	0.00
2	1,4-Dioxane	0.400	0.393	1.8	104	0.00
3	n-Nitrosodimethylamine	0.400	0.322	19.5	96	0.00
4 S	2-Fluorophenol	0.400	0.335	16.3	87	0.00
5 S	Phenol-d6	0.400	0.329	17.8	95	0.00
6	bis(2-Chloroethyl)ether	0.400	0.330	17.5	93	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	109	0.01
8 S	Nitrobenzene-d5	0.400	0.341	14.8	104	0.01
9	Naphthalene	0.400	0.404	-1.0	107	0.00
10	Hexachlorobutadiene	0.400	0.420	-5.0	110	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	108	0.00
12	2-Methylnaphthalene	0.400	0.388	3.0	107	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	104	0.00
14 S	2,4,6-Tribromophenol	0.400	0.324	19.0	97	0.00
15 S	2-Fluorobiphenyl	0.400	0.411	-2.7	105	0.00
16	Acenaphthylene	0.400	0.321	19.8	92	0.00
17	Acenaphthene	0.400	0.392	2.0	105	0.00
18	Fluorene	0.400	0.391	2.3	103	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	96	0.00
20	4-Bromophenyl-phenylether	0.400	0.353	11.8	89	0.00
21	Hexachlorobenzene	0.400	0.404	-1.0	96	-0.01
22	Pentachlorophenol	0.400	0.373	6.8	55	0.00
23	Phenanthrene	0.400	0.381	4.8	93	0.00
24	Anthracene	0.400	0.352	12.0	90	0.00
25 SURR	Fluoranthene-d10	0.400	0.351	12.3	87	0.00
26	Fluoranthene	0.400	0.345	13.8	89	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	90	0.00
28	Pvrene	0.400	0.386	3.5	85	0.00
29 S	Terphenyl-d14	0.400	0.409	-2.2	91	0.00
30	Benzo(a)anthracene	0.400	0.370	7.5	87	0.00
31	Chrysene	0.400	0.455	-13.7	100	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.328	18.0	85	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.395	1.3	91	0.00
34 I	Perylene-d12	0.400	0.400	0.0	92	0.00
35	Benzo(b)fluoranthene	0.400	0.383	4.3	94	0.00
36	Benzo(k)fluoranthene	0.400	0.412	-3.0	95	0.00
37 C	Benzo(a)pyrene	0.400	0.366	8.5	87	0.00
38	Dibenzo(a,h)anthracene	0.400	0.393	1.8	98	-0.01
39	Benzo(a,h,i)perylene	0.400	0.410	-2.5	92	-0.01

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

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