

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE020720\
 Data File : BE101135.D
 Acq On : 7 Feb 2020 13:12
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Feb 07 15:39:19 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	49	-0.02
2	1,4-Dioxane	0.400	0.412	-3.0	51	-0.02
3	n-Nitrosodimethylamine	0.400	0.387	3.3	55	-0.02
4 S	2-Fluorophenol	0.400	0.463	-15.8	57	-0.01
5 S	Phenol-d6	0.400	0.450	-12.5	61	0.01
6	bis(2-Chloroethyl)ether	0.400	0.283	29.3#	38	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	51	0.00
8 S	Nitrobenzene-d5	0.400	0.371	7.3	53	0.01
9	Naphthalene	0.400	0.396	1.0	50	0.00
10	Hexachlorobutadiene	0.400	0.436	-9.0	54	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	51	-0.01
12	2-Methylnaphthalene	0.400	0.366	8.5	48	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	58	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.183	54.3#	31	0.00
15 S	2-Fluorobiphenyl	0.400	0.359	10.3	51	-0.01
16	Acenaphthylene	0.400	0.383	4.3	61	-0.01
17	Acenaphthene	0.400	0.370	7.5	55	-0.03
18	Fluorene	0.400	0.379	5.3	56	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	64	-0.01
20	4-Bromophenyl-phenylether	0.400	0.403	-0.8	67	-0.01
21	Hexachlorobenzene	0.400	0.393	1.8	62	-0.01
22	Pentachlorophenol	0.400	0.251	37.3#	0	-0.04
23	Phenanthrene	0.400	0.351	12.3	57	-0.01
24	Anthracene	0.400	0.403	-0.8	68	0.00
25 SURR	Fluoranthene-d10	0.400	0.423	-5.7	70	-0.02
26	Fluoranthene	0.400	0.409	-2.2	69	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	81	-0.01
28	Pvrene	0.400	0.356	11.0	70	-0.01
29 S	Terphenyl-d14	0.400	0.362	9.5	72	-0.01
30	Benzo(a)anthracene	0.400	0.394	1.5	83	-0.01
31	Chrysene	0.400	0.392	2.0	77	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.540	-35.0#	125	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.435	-8.7	90	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	91	-0.01
35	Benzo(b)fluoranthene	0.400	0.331	17.3	81	-0.01
36	Benzo(k)fluoranthene	0.400	0.371	7.3	85	-0.01
37 C	Benzo(a)pyrene	0.400	0.388	3.0	92	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.400	0.0	99	-0.01
39	Benzo(a,h,i)perylene	0.400	0.357	10.8	79	-0.01

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

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