

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060519\
 Data File : BE099812.D
 Acq On : 5 Jun 2019 20:15
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc : MDL-SIM-0.1PPM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 06 01:41:07 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 05 14:35:58 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	82	0.00
2	1,4-Dioxane	0.400	0.355	11.3	85	0.00
3	n-Nitrosodimethylamine	0.400	0.263	34.3#	66	0.02
4 S	2-Fluorophenol	0.400	0.326	18.5	78	0.00
5 S	Phenol-d6	0.400	0.328	18.0	75	0.00
6	bis(2-Chloroethyl)ether	0.400	0.361	9.8	87	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	81	0.00
8 S	Nitrobenzene-d5	0.400	0.329	17.8	76	0.01
9	Naphthalene	0.400	0.363	9.3	82	0.00
10	Hexachlorobutadiene	0.400	0.363	9.3	82	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.354	11.5	80	0.01
12	2-Methylnaphthalene	0.400	0.337	15.8	78	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	76	0.01
14 S	2,4,6-Tribromophenol	0.400	0.367	8.3	68	0.00
15 S	2-Fluorobiphenyl	0.400	0.361	9.8	80	0.00
16	Acenaphthylene	0.400	0.368	8.0	79	0.00
17	Acenaphthene	0.400	0.356	11.0	77	0.00
18	Fluorene	0.400	0.370	7.5	77	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	75	0.00
20	4-Bromophenyl-phenylether	0.400	0.314	21.5	72	0.01
21	Hexachlorobenzene	0.400	0.343	14.2	75	0.00
22	Pentachlorophenol	0.400	0.363	9.3	40	0.03
23	Phenanthrene	0.400	0.347	13.3	74	0.00
24	Anthracene	0.400	0.318	20.5	72	0.00
25 SURR	Fluoranthene-d10	0.400	0.362	9.5	78	0.00
26	Fluoranthene	0.400	0.368	8.0	79	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	85	0.00
28	Pvrene	0.400	0.368	8.0	79	0.00
29 S	Terphenyl-d14	0.400	0.361	9.8	79	0.00
30	Benzo(a)anthracene	0.400	0.396	1.0	88	0.00
31	Chrysene	0.400	0.382	4.5	82	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.328	18.0	72	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.423	-5.7	92	0.00
34 I	Perylene-d12	0.400	0.400	0.0	90	0.00
35	Benzo(b)fluoranthene	0.400	0.365	8.8	92	0.00
36	Benzo(k)fluoranthene	0.400	0.347	13.3	87	0.00
37 C	Benzo(a)pyrene	0.400	0.365	8.8	91	0.00
38	Dibenzo(a,h)anthracene	0.400	0.351	12.3	90	0.00
39	Benzo(a,h,i)perylene	0.400	0.356	11.0	91	0.00

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

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