

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061719\
 Data File : BE100074.D
 Acq On : 17 Jun 2019 23:00
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Jun 18 03:59:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 17:23:36 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	88	0.00
2	1,4-Dioxane	0.400	0.365	8.8	81	0.00
3	n-Nitrosodimethylamine	0.400	0.569	-42.2#	130	0.01
4 S	2-Fluorophenol	0.400	0.346	13.5	81	0.00
5 S	Phenol-d6	0.400	0.380	5.0	88	0.00
6	bis(2-Chloroethyl)ether	0.400	0.433	-8.2	94	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	92	0.00
8 S	Nitrobenzene-d5	0.400	0.363	9.3	92	0.00
9	Naphthalene	0.400	0.407	-1.7	93	0.00
10	Hexachlorobutadiene	0.400	0.426	-6.5	97	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	94	0.00
12	2-Methylnaphthalene	0.400	0.393	1.8	93	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	89	0.00
14 S	2,4,6-Tribromophenol	0.400	0.346	13.5	89	0.00
15 S	2-Fluorobiphenyl	0.400	0.406	-1.5	89	0.00
16	Acenaphthylene	0.400	0.420	-5.0	95	0.00
17	Acenaphthene	0.400	0.411	-2.7	91	-0.02
18	Fluorene	0.400	0.424	-6.0	95	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	92	0.00
20	4-Bromophenyl-phenylether	0.400	0.392	2.0	93	0.00
21	Hexachlorobenzene	0.400	0.396	1.0	93	0.00
22	Pentachlorophenol	0.400	0.363	9.3	71	0.00
23	Phenanthrene	0.400	0.388	3.0	93	0.00
24	Anthracene	0.400	0.375	6.3	93	0.00
25 SURR	Fluoranthene-d10	0.400	0.369	7.8	89	0.00
26	Fluoranthene	0.400	0.381	4.8	91	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	87	0.00
28	Pvrene	0.400	0.451	-12.7	89	0.00
29 S	Terphenyl-d14	0.400	0.428	-7.0	86	0.00
30	Benzo(a)anthracene	0.400	0.421	-5.2	86	0.00
31	Chrysene	0.400	0.455	-13.7	89	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.328	18.0	71	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.469	-17.2	92	0.00
34 I	Perylene-d12	0.400	0.400	0.0	88	0.00
35	Benzo(b)fluoranthene	0.400	0.417	-4.2	90	0.00
36	Benzo(k)fluoranthene	0.400	0.426	-6.5	91	0.00
37 C	Benzo(a)pyrene	0.400	0.426	-6.5	93	0.00
38	Dibenzo(a,h)anthracene	0.400	0.417	-4.2	91	0.01
39	Benzo(a,h,i)perylene	0.400	0.417	-4.2	90	0.00

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