

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE060719\
 Data File : BE099871.D
 Acq On : 8 Jun 2019 2:17
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jun 08 04:35:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE060719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 07 15:25:21 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	104	0.00
2	1,4-Dioxane	0.400	0.414	-3.5	121	0.00
3	n-Nitrosodimethylamine	0.400	0.344	14.0	96	0.00
4 S	2-Fluorophenol	0.400	0.339	15.3	93	0.00
5 S	Phenol-d6	0.400	0.368	8.0	96	0.00
6	bis(2-Chloroethyl)ether	0.400	0.409	-2.2	116	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	113	-0.01
8 S	Nitrobenzene-d5	0.400	0.372	7.0	109	0.00
9	Naphthalene	0.400	0.398	0.5	114	0.00
10	Hexachlorobutadiene	0.400	0.408	-2.0	116	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.379	5.3	107	0.00
12	2-Methylnaphthalene	0.400	0.373	6.8	106	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.400	0.278	30.5#	67	-0.01
15 S	2-Fluorobiphenyl	0.400	0.413	-3.2	111	-0.01
16	Acenaphthylene	0.400	0.396	1.0	104	0.00
17	Acenaphthene	0.400	0.399	0.3	105	-0.01
18	Fluorene	0.400	0.383	4.3	98	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	96	-0.01
20	4-Bromophenyl-phenylether	0.400	0.400	0.0	101	0.00
21	Hexachlorobenzene	0.400	0.397	0.8	97	0.00
22	Pentachlorophenol	0.400	0.699	-74.7#	709	-0.04
23	Phenanthrene	0.400	0.398	0.5	95	-0.01
24	Anthracene	0.400	0.360	10.0	87	-0.01
25 SURR	Fluoranthene-d10	0.400	0.388	3.0	91	0.00
26	Fluoranthene	0.400	0.399	0.3	93	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	104	-0.01
28	Pvrene	0.400	0.400	0.0	91	0.00
29 S	Terphenyl-d14	0.400	0.390	2.5	91	0.00
30	Benzo(a)anthracene	0.400	0.414	-3.5	100	0.00
31	Chrysene	0.400	0.439	-9.7	104	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.348	13.0	84	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.442	-10.5	107	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	105	-0.01
35	Benzo(b)fluoranthene	0.400	0.399	0.3	107	-0.01
36	Benzo(k)fluoranthene	0.400	0.409	-2.2	106	-0.01
37 C	Benzo(a)pyrene	0.400	0.401	-0.3	106	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.400	0.0	106	-0.01
39	Benzo(a,h,i)perylene	0.400	0.407	-1.7	107	-0.02

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

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