

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE072318\
 Data File : BE097030.D
 Acq On : 23 Jul 2018 10:16
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 23 11:12:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE071618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 11:34:20 2018
 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	63	0.00
2	1,4-Dioxane	0.400	0.464	-16.0	76	0.00
3	n-Nitrosodimethylamine	0.400	0.419	-4.7	71	0.00
4 S	2-Fluorophenol	0.400	0.356	11.0	64	0.00
5 S	Phenol-d6	0.400	0.326	18.5	57	0.00
6	bis(2-Chloroethyl)ether	0.400	0.376	6.0	62	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	53	0.00
8 S	Nitrobenzene-d5	0.400	0.410	-2.5	59	0.00
9	Naphthalene	0.400	0.395	1.3	55	0.00
10	Hexachlorobutadiene	0.400	0.471	-17.7	66	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.378	5.5	53	0.00
12	2-Methylnaphthalene	0.400	0.375	6.3	52	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	57	0.00
14 S	2,4,6-Tribromophenol	0.400	0.385	3.8	57	0.00
15 S	2-Fluorobiphenyl	0.400	0.385	3.8	56	0.00
16	Acenaphthylene	0.400	0.377	5.8	56	0.00
17	Acenaphthene	0.400	0.363	9.3	54	0.00
18	Fluorene	0.400	0.398	0.5	60	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	70	0.00
20	4-Bromophenyl-phenylether	0.400	0.378	5.5	71	0.00
21	Hexachlorobenzene	0.400	0.388	3.0	71	0.00
22	Pentachlorophenol	0.400	0.591	-47.7#	112	0.00
23	Phenanthrene	0.400	0.384	4.0	71	0.00
24	Anthracene	0.400	0.372	7.0	70	0.00
25 SURR	Fluoranthene-d10	0.400	0.435	-8.7	83	0.00
26	Fluoranthene	0.400	0.411	-2.7	77	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	83	0.00
28	Pvrene	0.400	0.363	9.3	78	0.00
29 S	Terphenyl-d14	0.400	0.387	3.3	84	0.00
30	Benzo(a)anthracene	0.400	0.377	5.8	86	0.00
31	Chrysene	0.400	0.382	4.5	82	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.266	33.5#	64	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.402	-0.5	94	0.00
34 I	Perylene-d12	0.400	0.400	0.0	82	0.00
35	Benzo(b)fluoranthene	0.400	0.341	14.8	75	0.00
36	Benzo(k)fluoranthene	0.400	0.358	10.5	77	0.00
37 C	Benzo(a)pyrene	0.400	0.372	7.0	78	0.00
38	Dibenzo(a,h)anthracene	0.400	0.431	-7.7	97	0.00
39	Benzo(a,h,i)perylene	0.400	0.422	-5.5	96	0.00

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

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