

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE073118\
 Data File : BE097123.D
 Acq On : 31 Jul 2018 15:46
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 31 16:51:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE072418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 31 10:49:28 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	144	0.00
2	1,4-Dioxane	0.400	0.365	8.8	135	0.00
3	n-Nitrosodimethylamine	0.400	0.393	1.8	152	0.00
4 S	2-Fluorophenol	0.400	0.411	-2.7	157	0.00
5 S	Phenol-d6	0.400	0.433	-8.2	168	0.00
6	bis(2-Chloroethyl)ether	0.400	0.392	2.0	152	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	148	0.00
8 S	Nitrobenzene-d5	0.400	0.409	-2.2	160	0.00
9	Naphthalene	0.400	0.389	2.8	151	0.00
10	Hexachlorobutadiene	0.400	0.387	3.3	149	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.390	2.5	151	0.00
12	2-Methylnaphthalene	0.400	0.394	1.5	153	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	151	0.00
14 S	2,4,6-Tribromophenol	0.400	0.449	-12.2	188	0.00
15 S	2-Fluorobiphenyl	0.400	0.390	2.5	151	0.00
16	Acenaphthylene	0.400	0.390	2.5	152	0.00
17	Acenaphthene	0.400	0.395	1.3	153	0.00
18	Fluorene	0.400	0.400	0.0	158	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	154	0.00
20	4-Bromophenyl-phenylether	0.400	0.390	2.5	157	0.00
21	Hexachlorobenzene	0.400	0.370	7.5	145	0.00
22	Pentachlorophenol	0.400	0.640	-60.0#	286	0.01
23	Phenanthrene	0.400	0.391	2.3	158	0.00
24	Anthracene	0.400	0.398	0.5	165	0.01
25 SURR	Fluoranthene-d10	0.400	0.397	0.8	167	0.00
26	Fluoranthene	0.400	0.396	1.0	163	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	170	0.00
28	Pvrene	0.400	0.388	3.0	173	0.00
29 S	Terphenyl-d14	0.400	0.388	3.0	175	0.00
30	Benzo(a)anthracene	0.400	0.400	0.0	189	0.00
31	Chrysene	0.400	0.402	-0.5	176	0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.673	-68.3#	306	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.476	-19.0	217	0.00
34 I	Perylene-d12	0.400	0.400	0.0	204	0.00
35	Benzo(b)fluoranthene	0.400	0.372	7.0	201	0.00
36	Benzo(k)fluoranthene	0.400	0.364	9.0	200	0.00
37 C	Benzo(a)pyrene	0.400	0.393	1.8	213	0.00
38	Dibenzo(a,h)anthracene	0.400	0.402	-0.5	217	0.00
39	Benzo(a,h,i)perylene	0.400	0.389	2.8	209	0.00

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

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