

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100518\
 Data File : BE097759.D
 Acq On : 5 Oct 2018 17:40
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 06 04:19:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 03 12:17:19 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	60	-0.01
2	1,4-Dioxane	0.400	0.373	6.8	60	0.00
3	n-Nitrosodimethylamine	0.400	0.319	20.3	51	0.00
4 S	2-Fluorophenol	0.400	0.388	3.0	61	0.00
5 S	Phenol-d6	0.400	0.367	8.3	56	0.00
6	bis(2-Chloroethyl)ether	0.400	0.382	4.5	59	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	61	0.00
8 S	Nitrobenzene-d5	0.400	0.642	-60.5#	105	0.00
9	Naphthalene	0.400	0.386	3.5	60	0.00
10	Hexachlorobutadiene	0.400	0.390	2.5	62	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.385	3.8	60	0.00
12	2-Methylnaphthalene	0.400	0.376	6.0	60	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	63	0.00
14 S	2,4,6-Tribromophenol	0.400	0.450	-12.5	88	-0.01
15 S	2-Fluorobiphenyl	0.400	0.368	8.0	62	-0.02
16	Acenaphthylene	0.400	0.363	9.3	60	0.00
17	Acenaphthene	0.400	0.381	4.8	63	0.00
18	Fluorene	0.400	0.379	5.3	63	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	62	0.00
20	4-Bromophenyl-phenylether	0.400	0.373	6.8	62	-0.01
21	Hexachlorobenzene	0.400	0.373	6.8	61	0.00
22	Pentachlorophenol	0.400	0.441	-10.2	78	-0.01
23	Phenanthrene	0.400	0.377	5.8	62	0.00
24	Anthracene	0.400	0.357	10.8	61	0.00
25 SURR	Fluoranthene-d10	0.400	0.369	7.8	62	0.00
26	Fluoranthene	0.400	0.370	7.5	62	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	58	0.00
28	Pvrene	0.400	0.387	3.3	62	0.00
29 S	Terphenyl-d14	0.400	0.391	2.3	63	0.00
30	Benzo(a)anthracene	0.400	0.364	9.0	59	-0.01
31	Chrysene	0.400	0.406	-1.5	63	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.392	2.0	65	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.343	14.2	54	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	59	0.00
35	Benzo(b)fluoranthene	0.400	0.361	9.8	59	0.00
36	Benzo(k)fluoranthene	0.400	0.373	6.8	61	0.00
37 C	Benzo(a)pyrene	0.400	0.350	12.5	57	0.00
38	Dibenzo(a,h)anthracene	0.400	0.337	15.8	55	-0.02
39	Benzo(a,h,i)perylene	0.400	0.359	10.3	57	-0.02

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

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