

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE081018\
 Data File : BE097255.D
 Acq On : 10 Aug 2018 9:01
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 10 15:56:29 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	85	0.00
2	1,4-Dioxane	0.400	0.329	17.8	72	0.00
3	n-Nitrosodimethylamine	0.400	0.308	23.0	70	0.00
4 S	2-Fluorophenol	0.400	0.414	-3.5	93	0.00
5 S	Phenol-d6	0.400	0.432	-8.0	98	0.00
6	bis(2-Chloroethyl)ether	0.400	0.366	8.5	84	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	88	0.00
8 S	Nitrobenzene-d5	0.400	0.413	-3.2	96	0.00
9	Naphthalene	0.400	0.385	3.8	89	-0.01
10	Hexachlorobutadiene	0.400	0.397	0.8	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.367	8.3	85	0.00
12	2-Methylnaphthalene	0.400	0.372	7.0	86	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	80	0.00
14 S	2,4,6-Tribromophenol	0.400	0.386	3.5	85	0.00
15 S	2-Fluorobiphenyl	0.400	0.395	1.3	81	0.00
16	Acenaphthylene	0.400	0.386	3.5	79	0.00
17	Acenaphthene	0.400	0.380	5.0	78	-0.01
18	Fluorene	0.400	0.350	12.5	73	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	65	0.00
20	4-Bromophenyl-phenylether	0.400	0.414	-3.5	70	0.00
21	Hexachlorobenzene	0.400	0.436	-9.0	73	0.00
22	Pentachlorophenol	0.400	0.397	0.8	69	0.01
23	Phenanthrene	0.400	0.385	3.8	66	-0.01
24	Anthracene	0.400	0.390	2.5	68	0.00
25 SURR	Fluoranthene-d10	0.400	0.393	1.8	70	0.00
26	Fluoranthene	0.400	0.392	2.0	68	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	69	-0.01
28	Pyrene	0.400	0.390	2.5	71	0.00
29 S	Terphenyl-d14	0.400	0.375	6.3	69	0.00
30	Benzo(a)anthracene	0.400	0.393	1.8	76	0.00
31	Chrysene	0.400	0.385	3.8	69	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.567	-41.7#	103	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.508	-27.0#	94	0.00
34 I	Perylene-d12	0.400	0.400	0.0	85	0.00
35	Benzo(b)fluoranthene	0.400	0.361	9.8	81	0.00
36	Benzo(k)fluoranthene	0.400	0.371	7.3	85	0.00
37 C	Benzo(a)pyrene	0.400	0.396	1.0	89	0.00
38	Dibenzo(a,h)anthracene	0.400	0.420	-5.0	94	0.00
39	Benzo(g,h,i)perylene	0.400	0.390	2.5	87	0.01

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