

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121418\
 Data File : BE098452.D
 Acq On : 15 Dec 2018 1:14
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 17 17:03:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 14:59:55 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	88	-0.01
2	1,4-Dioxane	0.400	0.297	25.8#	66	0.00
3	n-Nitrosodimethylamine	0.400	0.362	9.5	86	0.00
4 S	2-Fluorophenol	0.400	0.486	-21.5	110	0.00
5 S	Phenol-d6	0.400	0.508	-27.0#	111	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.363	9.3	83	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	88	-0.01
8 S	Nitrobenzene-d5	0.400	0.406	-1.5	98	0.00
9	Naphthalene	0.400	0.389	2.8	88	0.00
10	Hexachlorobutadiene	0.400	0.401	-0.3	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.376	6.0	84	-0.01
12	2-Methylnaphthalene	0.400	0.381	4.8	86	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	87	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.505	-26.2#	116	0.00
15 S	2-Fluorobiphenyl	0.400	0.361	9.8	87	0.00
16	Acenaphthylene	0.400	0.389	2.8	88	0.00
17	Acenaphthene	0.400	0.380	5.0	84	0.00
18	Fluorene	0.400	0.390	2.5	86	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	85	-0.01
20	4-Bromophenyl-phenylether	0.400	0.365	8.8	81	0.00
21	Hexachlorobenzene	0.400	0.374	6.5	81	-0.01
22	Pentachlorophenol	0.400	0.693	-73.2#	166	-0.01
23	Phenanthrene	0.400	0.379	5.3	82	0.00
24	Anthracene	0.400	0.360	10.0	77	0.00
25 SURR	Fluoranthene-d10	0.400	0.401	-0.3	82	0.00
26	Fluoranthene	0.400	0.402	-0.5	82	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	101	-0.01
28	Pvrene	0.400	0.344	14.0	85	0.00
29 S	Terphenyl-d14	0.400	0.322	19.5	82	0.00
30	Benzo(a)anthracene	0.400	0.364	9.0	94	-0.01
31	Chrysene	0.400	0.379	5.3	96	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.398	0.5	101	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.393	1.8	104	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	107	-0.01
35	Benzo(b)fluoranthene	0.400	0.387	3.3	107	-0.01
36	Benzo(k)fluoranthene	0.400	0.378	5.5	105	-0.01
37 C	Benzo(a)pyrene	0.400	0.388	3.0	106	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.380	5.0	104	-0.02
39	Benzo(a,h,i)perylene	0.400	0.378	5.5	104	-0.02

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

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