

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110818\
 Data File : BE098179.D
 Acq On : 8 Nov 2018 9:28
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 09 09:48:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:00:52 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	99	0.00
2	1,4-Dioxane	0.400	0.409	-2.2	102	0.00
3	n-Nitrosodimethylamine	0.400	0.395	1.3	105	0.00
4 S	2-Fluorophenol	0.400	0.399	0.3	97	0.00
5 S	Phenol-d6	0.400	0.390	2.5	97	0.00
6	bis(2-Chloroethyl)ether	0.400	0.401	-0.3	103	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	0.00
8 S	Nitrobenzene-d5	0.400	0.393	1.8	101	-0.01
9	Naphthalene	0.400	0.405	-1.3	103	0.00
10	Hexachlorobutadiene	0.400	0.392	2.0	101	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.400	0.0	100	0.00
12	2-Methylnaphthalene	0.400	0.409	-2.2	105	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	93	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.330	17.5	91	0.00
15 S	2-Fluorobiphenyl	0.400	0.400	0.0	102	-0.01
16	Acenaphthylene	0.400	0.406	-1.5	102	0.00
17	Acenaphthene	0.400	0.408	-2.0	99	0.00
18	Fluorene	0.400	0.406	-1.5	101	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	93	0.00
20	4-Bromophenyl-phenylether	0.400	0.410	-2.5	98	0.00
21	Hexachlorobenzene	0.400	0.422	-5.5	103	0.00
22	Pentachlorophenol	0.400	0.418	-4.5	100	0.00
23	Phenanthrene	0.400	0.418	-4.5	101	0.00
24	Anthracene	0.400	0.405	-1.3	100	-0.01
25 SURR	Fluoranthene-d10	0.400	0.403	-0.8	99	0.00
26	Fluoranthene	0.400	0.412	-3.0	102	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	88	0.00
28	Pyrene	0.400	0.428	-7.0	103	0.00
29 S	Terphenyl-d14	0.400	0.414	-3.5	100	0.00
30	Benzo(a)anthracene	0.400	0.413	-3.2	99	0.00
31	Chrysene	0.400	0.416	-4.0	100	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.260	35.0#	71	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.387	3.3	93	0.00
34 I	Perylene-d12	0.400	0.400	0.0	88	0.00
35	Benzo(b)fluoranthene	0.400	0.421	-5.2	98	0.00
36	Benzo(k)fluoranthene	0.400	0.420	-5.0	99	0.00
37 C	Benzo(a)pyrene	0.400	0.410	-2.5	96	0.00
38	Dibenzo(a,h)anthracene	0.400	0.397	0.8	92	0.00
39	Benzo(g,h,i)perylene	0.400	0.399	0.3	93	0.00

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

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