

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE113018\
 Data File : BE098290.D
 Acq On : 30 Nov 2018 17:24
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 30 17:57:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE112918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 29 13:50:43 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	120	0.01
2	1,4-Dioxane	0.400	0.377	5.8	119	0.00
3	n-Nitrosodimethylamine	0.400	0.316	21.0	96	0.00
4 S	2-Fluorophenol	0.400	0.379	5.3	115	0.01
5 S	Phenol-d6	0.400	0.387	3.3	120	0.01
6	bis(2-Chloroethyl)ether	0.400	0.380	5.0	115	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	124	0.01
8 S	Nitrobenzene-d5	0.400	0.374	6.5	119	0.01
9	Naphthalene	0.400	0.394	1.5	123	0.00
10	Hexachlorobutadiene	0.400	0.390	2.5	122	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.400	0.0	121	0.01
12	2-Methylnaphthalene	0.400	0.393	1.8	122	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	126	0.01
14 S	2,4,6-Tribromophenol	0.400	0.337	15.8	120	0.01
15 S	2-Fluorobiphenyl	0.400	0.382	4.5	129	0.01
16	Acenaphthylene	0.400	0.393	1.8	127	0.00
17	Acenaphthene	0.400	0.383	4.3	124	0.01
18	Fluorene	0.400	0.401	-0.3	127	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	134	0.00
20	4-Bromophenyl-phenylether	0.400	0.360	10.0	123	0.01
21	Hexachlorobenzene	0.400	0.375	6.3	126	0.01
22	Pentachlorophenol	0.400	0.298	25.5#	108	0.00
23	Phenanthrene	0.400	0.385	3.8	131	0.01
24	Anthracene	0.400	0.371	7.3	129	0.01
25 SURR	Fluoranthene-d10	0.400	0.373	6.8	128	0.00
26	Fluoranthene	0.400	0.376	6.0	131	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	121	0.00
28	Pvrene	0.400	0.465	-16.3	131	0.00
29 S	Terphenyl-d14	0.400	0.431	-7.7	125	0.00
30	Benzo(a)anthracene	0.400	0.378	5.5	118	0.00
31	Chrysene	0.400	0.400	0.0	123	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.351	12.3	108	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.386	3.5	125	0.00
34 I	Perylene-d12	0.400	0.400	0.0	124	0.00
35	Benzo(b)fluoranthene	0.400	0.381	4.8	120	0.00
36	Benzo(k)fluoranthene	0.400	0.405	-1.3	130	0.00
37 C	Benzo(a)pyrene	0.400	0.394	1.5	124	0.00
38	Dibenzo(a,h)anthracene	0.400	0.396	1.0	125	0.00
39	Benzo(a,h,i)perylene	0.400	0.391	2.3	125	0.00

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

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