

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE112918\
 Data File : BE098281.D
 Acq On : 29 Nov 2018 19:32
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 30 00:30:43 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	128	0.00
2	1,4-Dioxane	0.400	0.328	18.0	112	0.00
3	n-Nitrosodimethylamine	0.400	0.359	10.3	116	0.00
4 S	2-Fluorophenol	0.400	0.397	0.8	128	0.01
5 S	Phenol-d6	0.400	0.409	-2.2	135	0.00
6	bis(2-Chloroethyl)ether	0.400	0.385	3.8	125	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	128	0.00
8 S	Nitrobenzene-d5	0.400	0.386	3.5	127	0.00
9	Naphthalene	0.400	0.397	0.8	128	-0.01
10	Hexachlorobutadiene	0.400	0.389	2.8	126	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.409	-2.2	127	0.00
12	2-Methylnaphthalene	0.400	0.396	1.0	127	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	130	0.00
14 S	2,4,6-Tribromophenol	0.400	0.368	8.0	136	0.00
15 S	2-Fluorobiphenyl	0.400	0.376	6.0	131	0.00
16	Acenaphthylene	0.400	0.387	3.3	129	-0.01
17	Acenaphthene	0.400	0.389	2.8	130	0.00
18	Fluorene	0.400	0.400	0.0	131	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	129	0.00
20	4-Bromophenyl-phenylether	0.400	0.387	3.3	127	0.00
21	Hexachlorobenzene	0.400	0.389	2.8	125	0.00
22	Pentachlorophenol	0.400	0.372	7.0	141	0.00
23	Phenanthrene	0.400	0.391	2.3	128	0.00
24	Anthracene	0.400	0.389	2.8	129	0.00
25 SURR	Fluoranthene-d10	0.400	0.372	7.0	123	0.00
26	Fluoranthene	0.400	0.369	7.8	123	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	120	0.00
28	Pvrene	0.400	0.443	-10.7	123	0.00
29 S	Terphenyl-d14	0.400	0.419	-4.7	121	0.00
30	Benzo(a)anthracene	0.400	0.383	4.3	118	0.00
31	Chrysene	0.400	0.386	3.5	118	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.396	1.0	121	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.376	6.0	121	0.00
34 I	Perylene-d12	0.400	0.400	0.0	120	0.00
35	Benzo(b)fluoranthene	0.400	0.391	2.3	119	0.00
36	Benzo(k)fluoranthene	0.400	0.394	1.5	122	0.00
37 C	Benzo(a)pyrene	0.400	0.396	1.0	121	0.00
38	Dibenzo(a,h)anthracene	0.400	0.392	2.0	120	0.00
39	Benzo(a,h,i)perylene	0.400	0.390	2.5	120	0.00

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

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