

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121218\
 Data File : BE098379.D
 Acq On : 13 Dec 2018 2:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 13 07:25:10 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	77	0.00
2	1,4-Dioxane	0.400	0.371	7.3	73	0.00
3	n-Nitrosodimethylamine	0.400	0.254	36.5#	53	0.00
4 S	2-Fluorophenol	0.400	0.366	8.5	73	0.01
5 S	Phenol-d6	0.400	0.364	9.0	70	0.00
6	bis(2-Chloroethyl)ether	0.400	0.323	19.3	64	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	76	-0.01
8 S	Nitrobenzene-d5	0.400	0.382	4.5	79	0.01
9	Naphthalene	0.400	0.392	2.0	76	0.00
10	Hexachlorobutadiene	0.400	0.419	-4.7	82	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.387	3.3	74	0.00
12	2-Methylnaphthalene	0.400	0.372	7.0	72	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	76	0.00
14 S	2,4,6-Tribromophenol	0.400	0.395	1.3	79	0.00
15 S	2-Fluorobiphenyl	0.400	0.361	9.8	76	0.00
16	Acenaphthylene	0.400	0.389	2.8	77	0.00
17	Acenaphthene	0.400	0.384	4.0	74	0.00
18	Fluorene	0.400	0.382	4.5	73	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	74	-0.01
20	4-Bromophenyl-phenylether	0.400	0.354	11.5	68	0.00
21	Hexachlorobenzene	0.400	0.387	3.3	73	0.00
22	Pentachlorophenol	0.400	0.307	23.3	56	0.00
23	Phenanthrene	0.400	0.373	6.8	70	0.00
24	Anthracene	0.400	0.343	14.2	64	0.00
25 SURR	Fluoranthene-d10	0.400	0.390	2.5	70	0.00
26	Fluoranthene	0.400	0.398	0.5	71	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	84	0.00
28	Pvrene	0.400	0.347	13.3	71	0.00
29 S	Terphenyl-d14	0.400	0.324	19.0	69	0.00
30	Benzo(a)anthracene	0.400	0.368	8.0	79	-0.01
31	Chrysene	0.400	0.377	5.8	80	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.324	19.0	69	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.390	2.5	87	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	90	0.00
35	Benzo(b)fluoranthene	0.400	0.392	2.0	91	0.00
36	Benzo(k)fluoranthene	0.400	0.398	0.5	93	-0.01
37 C	Benzo(a)pyrene	0.400	0.391	2.3	90	0.00
38	Dibenzo(a,h)anthracene	0.400	0.374	6.5	86	0.00
39	Benzo(a,h,i)perylene	0.400	0.378	5.5	87	0.00

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

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