

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121818\
 Data File : BE098520.D
 Acq On : 18 Dec 2018 19:39
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 19 03:25:31 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	73	-0.01
2	1,4-Dioxane	0.400	0.373	6.8	69	-0.01
3	n-Nitrosodimethylamine	0.400	0.290	27.5#	57	0.00
4 S	2-Fluorophenol	0.400	0.433	-8.2	81	0.00
5 S	Phenol-d6	0.400	0.507	-26.7#	91	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.386	3.5	73	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	85	-0.01
8 S	Nitrobenzene-d5	0.400	0.341	14.8	79	0.00
9	Naphthalene	0.400	0.397	0.8	87	-0.01
10	Hexachlorobutadiene	0.400	0.402	-0.5	88	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.411	-2.7	89	-0.01
12	2-Methylnaphthalene	0.400	0.399	0.3	87	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	82	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.368	8.0	79	-0.01
15 S	2-Fluorobiphenyl	0.400	0.396	1.0	89	-0.01
16	Acenaphthylene	0.400	0.354	11.5	75	-0.01
17	Acenaphthene	0.400	0.386	3.5	80	-0.01
18	Fluorene	0.400	0.368	8.0	76	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	69	-0.02
20	4-Bromophenyl-phenylether	0.400	0.391	2.3	70	-0.01
21	Hexachlorobenzene	0.400	0.422	-5.5	74	-0.01
22	Pentachlorophenol	0.400	0.408	-2.0	72	-0.01
23	Phenanthrene	0.400	0.384	4.0	67	-0.01
24	Anthracene	0.400	0.354	11.5	62	-0.01
25 SURR	Fluoranthene-d10	0.400	0.375	6.3	62	-0.01
26	Fluoranthene	0.400	0.370	7.5	61	-0.01
27 I	Chrysene-d12	0.400	0.400	0.0	74	-0.01
28	Pvrene	0.400	0.349	12.8	63	-0.01
29 S	Terphenyl-d14	0.400	0.340	15.0	64	-0.01
30	Benzo(a)anthracene	0.400	0.371	7.3	70	-0.01
31	Chrysene	0.400	0.380	5.0	71	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.330	17.5	61	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.285	28.8#	56	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	76	-0.03
35	Benzo(b)fluoranthene	0.400	0.389	2.8	77	-0.02
36	Benzo(k)fluoranthene	0.400	0.416	-4.0	83	-0.03
37 C	Benzo(a)pyrene	0.400	0.391	2.3	76	-0.03
38	Dibenzo(a,h)anthracene	0.400	0.310	22.5	61	-0.03
39	Benzo(a,h,i)perylene	0.400	0.283	29.3#	56	-0.03

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

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