

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE110118\
 Data File : BE098145.D
 Acq On : 1 Nov 2018 14:18
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 01 14:51:52 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	83	0.00
2	1,4-Dioxane	0.400	0.432	-8.0	90	0.00
3	n-Nitrosodimethylamine	0.400	0.420	-5.0	93	0.00
4 S	2-Fluorophenol	0.400	0.439	-9.7	89	0.00
5 S	Phenol-d6	0.400	0.430	-7.5	89	0.00
6	bis(2-Chloroethyl)ether	0.400	0.406	-1.5	87	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	87	0.00
8 S	Nitrobenzene-d5	0.400	0.395	1.3	88	0.00
9	Naphthalene	0.400	0.413	-3.2	91	0.00
10	Hexachlorobutadiene	0.400	0.420	-5.0	93	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	89	0.00
12	2-Methylnaphthalene	0.400	0.416	-4.0	93	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	84	0.00
14 S	2,4,6-Tribromophenol	0.400	0.361	9.8	90	0.00
15 S	2-Fluorobiphenyl	0.400	0.406	-1.5	94	0.00
16	Acenaphthylene	0.400	0.408	-2.0	92	0.00
17	Acenaphthene	0.400	0.413	-3.2	90	0.00
18	Fluorene	0.400	0.399	0.3	89	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	87	0.00
20	4-Bromophenyl-phenylether	0.400	0.413	-3.2	92	0.00
21	Hexachlorobenzene	0.400	0.416	-4.0	94	0.00
22	Pentachlorophenol	0.400	0.380	5.0	80	0.00
23	Phenanthrene	0.400	0.407	-1.7	92	0.00
24	Anthracene	0.400	0.409	-2.2	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.401	-0.3	92	0.00
26	Fluoranthene	0.400	0.400	0.0	92	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	81	0.00
28	Pvrene	0.400	0.420	-5.0	92	0.00
29 S	Terphenyl-d14	0.400	0.415	-3.7	92	0.00
30	Benzo(a)anthracene	0.400	0.415	-3.7	91	0.00
31	Chrysene	0.400	0.415	-3.7	91	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.483	-20.7	100	0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.411	-2.7	90	0.00
34 I	Perylene-d12	0.400	0.400	0.0	82	0.00
35	Benzo(b)fluoranthene	0.400	0.409	-2.2	89	0.00
36	Benzo(k)fluoranthene	0.400	0.420	-5.0	92	0.00
37 C	Benzo(a)pyrene	0.400	0.412	-3.0	89	0.00
38	Dibenzo(a,h)anthracene	0.400	0.413	-3.2	89	0.00
39	Benzo(a,h,i)perylene	0.400	0.416	-4.0	90	0.00

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

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