

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121318\
 Data File : BE098397.D
 Acq On : 13 Dec 2018 15:16
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 13 20:57:25 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	80	0.00
2	1,4-Dioxane	0.400	0.367	8.3	75	0.00
3	n-Nitrosodimethylamine	0.400	0.242	39.5#	52	0.01
4 S	2-Fluorophenol	0.400	0.379	5.3	78	0.01
5 S	Phenol-d6	0.400	0.392	2.0	78	0.00
6	bis(2-Chloroethyl)ether	0.400	0.323	19.3	67	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	79	-0.01
8 S	Nitrobenzene-d5	0.400	0.382	4.5	82	0.01
9	Naphthalene	0.400	0.389	2.8	79	0.00
10	Hexachlorobutadiene	0.400	0.406	-1.5	82	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.374	6.5	75	0.00
12	2-Methylnaphthalene	0.400	0.359	10.3	72	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	79	0.00
14 S	2,4,6-Tribromophenol	0.400	0.386	3.5	80	0.00
15 S	2-Fluorobiphenyl	0.400	0.363	9.3	79	0.00
16	Acenaphthylene	0.400	0.383	4.3	78	0.00
17	Acenaphthene	0.400	0.379	5.3	75	0.00
18	Fluorene	0.400	0.383	4.3	76	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	76	-0.01
20	4-Bromophenyl-phenylether	0.400	0.341	14.8	68	0.00
21	Hexachlorobenzene	0.400	0.383	4.3	74	0.00
22	Pentachlorophenol	0.400	0.318	20.5	60	0.00
23	Phenanthrene	0.400	0.378	5.5	73	0.00
24	Anthracene	0.400	0.340	15.0	66	0.00
25 SURR	Fluoranthene-d10	0.400	0.396	1.0	72	0.00
26	Fluoranthene	0.400	0.402	-0.5	73	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	88	0.00
28	Pvrene	0.400	0.345	13.8	74	0.00
29 S	Terphenyl-d14	0.400	0.316	21.0	70	0.00
30	Benzo(a)anthracene	0.400	0.361	9.8	81	-0.01
31	Chrysene	0.400	0.372	7.0	82	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.302	24.5	67	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.386	3.5	89	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	93	-0.01
35	Benzo(b)fluoranthene	0.400	0.380	5.0	91	0.00
36	Benzo(k)fluoranthene	0.400	0.405	-1.3	98	-0.01
37 C	Benzo(a)pyrene	0.400	0.375	6.3	89	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.369	7.8	88	-0.01
39	Benzo(a,h,i)perylene	0.400	0.373	6.8	89	0.00

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

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