

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE031119\
 Data File : BE098872.D
 Acq On : 11 Mar 2019 21:21
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 12 16:36:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE030619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 06 15:33:24 2019
 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	107	0.00
2	1,4-Dioxane	0.400	0.379	5.3	105	0.00
3	n-Nitrosodimethylamine	0.400	0.252	37.0#	66	0.02
4 S	2-Fluorophenol	0.400	0.346	13.5	89	0.01
5 S	Phenol-d6	0.400	0.383	4.3	103	0.00
6	bis(2-Chloroethyl)ether	0.400	0.354	11.5	98	0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	115	0.00
8 S	Nitrobenzene-d5	0.400	0.358	10.5	114	0.00
9	Naphthalene	0.400	0.390	2.5	115	0.00
10	Hexachlorobutadiene	0.400	0.377	5.8	112	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.392	2.0	115	0.00
12	2-Methylnaphthalene	0.400	0.387	3.3	114	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	122	0.00
14 S	2,4,6-Tribromophenol	0.400	0.315	21.3	97	0.00
15 S	2-Fluorobiphenyl	0.400	0.394	1.5	122	0.00
16	Acenaphthylene	0.400	0.375	6.3	120	0.00
17	Acenaphthene	0.400	0.372	7.0	114	-0.02
18	Fluorene	0.400	0.372	7.0	116	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
20	4-Bromophenyl-phenylether	0.400	0.373	6.8	104	0.00
21	Hexachlorobenzene	0.400	0.401	-0.3	116	0.00
22	Pentachlorophenol	0.400	0.407	-1.7	89	0.01
23	Phenanthrene	0.400	0.390	2.5	109	0.00
24	Anthracene	0.400	0.363	9.3	105	0.00
25 SURR	Fluoranthene-d10	0.400	0.363	9.3	103	0.00
26	Fluoranthene	0.400	0.364	9.0	103	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	-0.01
28	Pvrene	0.400	0.411	-2.7	102	0.00
29 S	Terphenyl-d14	0.400	0.392	2.0	98	0.00
30	Benzo(a)anthracene	0.400	0.393	1.8	100	0.00
31	Chrysene	0.400	0.383	4.3	97	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.313	21.8	85	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.404	-1.0	102	0.00
34 I	Perylene-d12	0.400	0.400	0.0	105	0.00
35	Benzo(b)fluoranthene	0.400	0.391	2.3	106	0.00
36	Benzo(k)fluoranthene	0.400	0.399	0.3	110	0.00
37 C	Benzo(a)pyrene	0.400	0.381	4.8	103	0.00
38	Dibenzo(a,h)anthracene	0.400	0.382	4.5	101	-0.01
39	Benzo(a,h,i)perylene	0.400	0.380	5.0	102	-0.01

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

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