

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121118\
 Data File : BE098360.D
 Acq On : 12 Dec 2018 13:38
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 12 14:10:39 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	135	0.00
2	1,4-Dioxane	0.400	0.399	0.3	137	0.00
3	n-Nitrosodimethylamine	0.400	0.356	11.0	129	0.00
4 S	2-Fluorophenol	0.400	0.410	-2.5	142	0.00
5 S	Phenol-d6	0.400	0.406	-1.5	136	0.00
6	bis(2-Chloroethyl)ether	0.400	0.357	10.8	124	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	128	0.01
8 S	Nitrobenzene-d5	0.400	0.445	-11.2	155	0.03
9	Naphthalene	0.400	0.390	2.5	128	0.01
10	Hexachlorobutadiene	0.400	0.402	-0.5	132	0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.376	6.0	121	0.01
12	2-Methylnaphthalene	0.400	0.374	6.5	122	0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	117	0.00
14 S	2,4,6-Tribromophenol	0.400	0.394	1.5	122	0.01
15 S	2-Fluorobiphenyl	0.400	0.389	2.8	125	0.01
16	Acenaphthylene	0.400	0.385	3.8	117	0.01
17	Acenaphthene	0.400	0.389	2.8	115	0.01
18	Fluorene	0.400	0.377	5.8	111	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	110	0.00
20	4-Bromophenyl-phenylether	0.400	0.383	4.3	110	0.01
21	Hexachlorobenzene	0.400	0.396	1.0	111	0.00
22	Pentachlorophenol	0.400	0.467	-16.8	135	0.00
23	Phenanthrene	0.400	0.389	2.8	108	0.00
24	Anthracene	0.400	0.369	7.8	103	0.01
25 SURR	Fluoranthene-d10	0.400	0.390	2.5	103	0.00
26	Fluoranthene	0.400	0.390	2.5	103	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	127	0.00
28	Pvrene	0.400	0.347	13.3	108	0.00
29 S	Terphenyl-d14	0.400	0.335	16.3	108	0.00
30	Benzo(a)anthracene	0.400	0.362	9.5	118	0.00
31	Chrysene	0.400	0.387	3.3	124	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.372	7.0	119	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.393	1.8	132	0.00
34 I	Perylene-d12	0.400	0.400	0.0	134	0.00
35	Benzo(b)fluoranthene	0.400	0.381	4.8	132	0.00
36	Benzo(k)fluoranthene	0.400	0.402	-0.5	140	0.00
37 C	Benzo(a)pyrene	0.400	0.387	3.3	132	0.00
38	Dibenzo(a,h)anthracene	0.400	0.387	3.3	133	0.00
39	Benzo(a,h,i)perylene	0.400	0.381	4.8	131	0.00

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

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