

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE121418\
 Data File : BE098454.D
 Acq On : 15 Dec 2018 3:41
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 15 04:15:17 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	103	-0.01
2	1,4-Dioxane	0.400	0.333	16.8	87	0.00
3	n-Nitrosodimethylamine	0.400	0.291	27.3#	81	0.00
4 S	2-Fluorophenol	0.400	0.429	-7.2	114	0.00
5 S	Phenol-d6	0.400	0.429	-7.2	110	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.333	16.8	88	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	99	-0.01
8 S	Nitrobenzene-d5	0.400	0.375	6.3	101	0.00
9	Naphthalene	0.400	0.395	1.3	100	0.00
10	Hexachlorobutadiene	0.400	0.407	-1.7	103	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.379	5.3	95	-0.01
12	2-Methylnaphthalene	0.400	0.374	6.5	94	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	97	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.461	-15.3	118	0.00
15 S	2-Fluorobiphenyl	0.400	0.374	6.5	99	0.00
16	Acenaphthylene	0.400	0.391	2.3	98	0.00
17	Acenaphthene	0.400	0.378	5.5	92	0.00
18	Fluorene	0.400	0.384	4.0	93	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	92	-0.01
20	4-Bromophenyl-phenylether	0.400	0.371	7.3	89	0.00
21	Hexachlorobenzene	0.400	0.389	2.8	91	0.00
22	Pentachlorophenol	0.400	0.553	-38.3#	138	0.00
23	Phenanthrene	0.400	0.376	6.0	88	0.00
24	Anthracene	0.400	0.356	11.0	83	0.00
25 SURR	Fluoranthene-d10	0.400	0.382	4.5	85	0.00
26	Fluoranthene	0.400	0.388	3.0	86	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	105	-0.01
28	Pvrene	0.400	0.346	13.5	89	0.00
29 S	Terphenyl-d14	0.400	0.316	21.0	84	0.00
30	Benzo(a)anthracene	0.400	0.363	9.3	97	-0.01
31	Chrysene	0.400	0.381	4.8	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.367	8.3	97	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.400	0.0	111	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	113	-0.02
35	Benzo(b)fluoranthene	0.400	0.387	3.3	113	-0.01
36	Benzo(k)fluoranthene	0.400	0.393	1.8	116	-0.01
37 C	Benzo(a)pyrene	0.400	0.388	3.0	112	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.378	5.5	110	-0.02
39	Benzo(a,h,i)perylene	0.400	0.377	5.8	110	-0.02

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

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