

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE110819\
 Data File : BE100711.D
 Acq On : 8 Nov 2019 10:49
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 08 12:40:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 16:49:33 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	41#	0.01
2	1,4-Dioxane	0.714	0.574	19.6	33#	0.00
3	n-Nitrosodimethylamine	0.755	0.671	11.1	39#	0.00
4 S	2-Fluorophenol	1.235	1.365	-10.5	47#	0.01
5 S	Phenol-d6	1.742	2.116	-21.5	52	0.02
6	bis(2-Chloroethyl)ether	1.430	1.373	4.0	42#	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	46#	0.01
8 S	Nitrobenzene-d5	0.248	0.361	-45.6#	77	0.01
9	Naphthalene	4.206	4.053	3.6	46#	0.01
10	Hexachlorobutadiene	0.183	0.183	0.0	48#	0.00
11 SURR	2-Methylnaphthalene-d10	0.631	0.636	-0.8	49#	0.01
12	2-Methylnaphthalene	0.717	0.712	0.7	48#	0.01
13 I	Acenaphthene-d10	1.000	1.000	0.0	50#	0.01
14 S	2,4,6-Tribromophenol	0.106	0.046	56.6#	24#	0.03
15 S	2-Fluorobiphenyl	1.457	1.488	-2.1	52	0.01
16	Acenaphthylene	1.721	1.636	4.9	49#	0.00
17	Acenaphthene	1.112	1.021	8.2	47#	0.01
18	Fluorene	1.477	1.392	5.8	48#	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	48#	0.00
20	4-Bromophenyl-phenylether	0.204	0.205	-0.5	50#	0.00
21	Hexachlorobenzene	0.194	0.195	-0.5	51	0.00
22	Pentachlorophenol	0.062	0.001	98.4#	1#	0.00
23	Phenanthrene	1.100	1.092	0.7	47#	0.00
24	Anthracene	1.021	1.009	1.2	49#	0.01
25 SURR	Fluoranthene-d10	5.162	4.868	5.7	47#	0.01
26	Fluoranthene	1.436	1.381	3.8	46#	0.01
27 I	Chrysene-d12	1.000	1.000	0.0	41#	0.01
28	Pvrene	1.118	1.244	-11.3	46#	0.00
29 S	Terphenyl-d14	0.875	1.036	-18.4	52	0.00
30	Benzo(a)anthracene	1.305	1.301	0.3	42#	0.01
31	Chrysene	1.304	1.255	3.8	40#	0.01
32	Bis(2-ethylhexyl)phthalate	2.729	3.299	-20.9	50#	0.00
33	Indeno(1,2,3-cd)pyrene	1.603	1.520	5.2	41#	0.04
34 I	Perylene-d12	1.000	1.000	0.0	38#	0.02
35	Benzo(b)fluoranthene	1.166	1.123	3.7	39#	0.02
36	Benzo(k)fluoranthene	1.200	1.138	5.2	38#	0.01
37 C	Benzo(a)pyrene	1.086	1.092	-0.6	41#	0.02
38	Dibenzo(a,h)anthracene	1.122	1.104	1.6	41#	0.04
39	Benzo(a,h,i)perylene	1.129	1.083	4.1	39#	0.04

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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