

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102319\
 Data File : BE100662.D
 Acq On : 23 Oct 2019 15:27
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 23 16:58:05 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	54	0.00
2	1,4-Dioxane	0.714	0.720	-0.8	55	0.00
3	n-Nitrosodimethylamine	0.755	0.743	1.6	58	0.00
4 S	2-Fluorophenol	1.235	1.324	-7.2	62	0.00
5 S	Phenol-d6	1.742	1.908	-9.5	63	0.00
6	bis(2-Chloroethyl)ether	1.430	1.386	3.1	56	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	57	0.00
8 S	Nitrobenzene-d5	0.248	0.349	-40.7#	91	0.00
9	Naphthalene	4.206	4.095	2.6	57	0.00
10	Hexachlorobutadiene	0.183	0.183	0.0	59	0.00
11 SURR	2-Methylnaphthalene-d10	0.631	0.625	1.0	58	0.00
12	2-Methylnaphthalene	0.717	0.713	0.6	59	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	59	0.00
14 S	2,4,6-Tribromophenol	0.106	0.171	-61.3#	109	0.00
15 S	2-Fluorobiphenyl	1.457	1.555	-6.7	65	0.00
16	Acenaphthylene	1.721	1.730	-0.5	62	0.00
17	Acenaphthene	1.112	1.086	2.3	60	0.00
18	Fluorene	1.477	1.454	1.6	60	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	64	0.00
20	4-Bromophenyl-phenylether	0.204	0.201	1.5	65	0.00
21	Hexachlorobenzene	0.194	0.191	1.5	67	0.00
22	Pentachlorophenol	0.062	0.101	-62.9#	117	0.00
23	Phenanthrene	1.100	1.122	-2.0	65	0.00
24	Anthracene	1.021	0.972	4.8	63	0.00
25 SURR	Fluoranthene-d10	5.162	5.346	-3.6	69	0.00
26	Fluoranthene	1.436	1.543	-7.5	69	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	71	0.00
28	Pvrene	1.118	1.117	0.1	71	0.00
29 S	Terphenyl-d14	0.875	0.948	-8.3	81	0.00
30	Benzo(a)anthracene	1.305	1.261	3.4	69	0.00
31	Chrysene	1.304	1.280	1.8	70	0.00
32	Bis(2-ethylhexyl)phthalate	2.729	3.096	-13.4	80	0.00
33	Indeno(1,2,3-cd)pyrene	1.603	1.482	7.5	68	0.00
34 I	Perylene-d12	1.000	1.000	0.0	64	0.00
35	Benzo(b)fluoranthene	1.166	1.122	3.8	67	0.00
36	Benzo(k)fluoranthene	1.200	1.146	4.5	66	0.00
37 C	Benzo(a)pyrene	1.086	1.033	4.9	66	0.00
38	Dibenzo(a,h)anthracene	1.122	1.042	7.1	66	0.00
39	Benzo(a,h,i)perylene	1.129	1.080	4.3	66	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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