

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE102819\
 Data File : BE100675.D
 Acq On : 28 Oct 2019 16:02
 Operator : JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 29 11:52:15 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	47#	0.00
2	1,4-Dioxane	0.714	0.856	-19.9	57	0.00
3	n-Nitrosodimethylamine	0.755	0.791	-4.8	54	0.00
4 S	2-Fluorophenol	1.235	1.275	-3.2	52	0.00
5 S	Phenol-d6	1.742	1.758	-0.9	51	0.00
6	bis(2-Chloroethyl)ether	1.430	1.633	-14.2	58	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	50#	0.00
8 S	Nitrobenzene-d5	0.248	0.321	-29.4#	74	0.00
9	Naphthalene	4.206	4.690	-11.5	57	0.00
10	Hexachlorobutadiene	0.183	0.210	-14.8	60	0.00
11 SURR	2-Methylnaphthalene-d10	0.631	0.683	-8.2	56	0.00
12	2-Methylnaphthalene	0.717	0.797	-11.2	58	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	50#	0.00
14 S	2,4,6-Tribromophenol	0.106	0.116	-9.4	62	0.00
15 S	2-Fluorobiphenyl	1.457	1.575	-8.1	55	0.00
16	Acenaphthylene	1.721	1.824	-6.0	55	0.00
17	Acenaphthene	1.112	1.217	-9.4	56	0.00
18	Fluorene	1.477	1.608	-8.9	56	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	51	-0.01
20	4-Bromophenyl-phenylether	0.204	0.224	-9.8	59	0.00
21	Hexachlorobenzene	0.194	0.216	-11.3	60	0.00
22	Pentachlorophenol	0.062	0.059	4.8	55	0.00
23	Phenanthrene	1.100	1.225	-11.4	57	-0.01
24	Anthracene	1.021	1.056	-3.4	55	0.00
25 SURR	Fluoranthene-d10	5.162	5.759	-11.6	60	0.00
26	Fluoranthene	1.436	1.630	-13.5	59	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	51	0.00
28	Pvrene	1.118	1.259	-12.6	57	0.00
29 S	Terphenyl-d14	0.875	0.969	-10.7	59	0.00
30	Benzo(a)anthracene	1.305	1.436	-10.0	56	0.00
31	Chrysene	1.304	1.491	-14.3	58	0.00
32	Bis(2-ethylhexyl)phthalate	2.729	2.458	9.9	46#	-0.01
33	Indeno(1,2,3-cd)pyrene	1.603	1.762	-9.9	58	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	49#	0.00
35	Benzo(b)fluoranthene	1.166	1.251	-7.3	57	0.00
36	Benzo(k)fluoranthene	1.200	1.301	-8.4	57	0.00
37 C	Benzo(a)pyrene	1.086	1.141	-5.1	56	0.00
38	Dibenzo(a,h)anthracene	1.122	1.209	-7.8	59	0.00
39	Benzo(a,h,i)perylene	1.129	1.250	-10.7	58	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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