

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE100518\
 Data File : BE097745.D
 Acq On : 5 Oct 2018 8:44
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 05 14:03:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE100318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 03 12:17:19 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	64	-0.01
2	1,4-Dioxane	0.797	0.723	9.3	62	0.00
3	n-Nitrosodimethylamine	0.725	0.593	18.2	55	0.00
4 S	2-Fluorophenol	0.946	0.904	4.4	64	0.00
5 S	Phenol-d6	1.450	1.345	7.2	61	0.00
6	bis(2-Chloroethyl)ether	1.501	1.401	6.7	62	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	64	0.00
8 S	Nitrobenzene-d5	0.130	0.203	-56.2#	107	0.00
9	Naphthalene	4.031	3.789	6.0	61	0.00
10	Hexachlorobutadiene	0.214	0.207	3.3	64	-0.01
11 SURR	2-Methylnaphthalene-d10	0.636	0.610	4.1	63	0.00
12	2-Methylnaphthalene	0.645	0.606	6.0	62	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	66	0.00
14 S	2,4,6-Tribromophenol	0.072	0.078	-8.3	88	-0.01
15 S	2-Fluorobiphenyl	1.761	1.595	9.4	64	-0.02
16	Acenaphthylene	1.771	1.564	11.7	61	0.00
17	Acenaphthene	1.135	1.037	8.6	64	0.00
18	Fluorene	1.507	1.400	7.1	65	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	64	0.00
20	4-Bromophenyl-phenylether	0.210	0.198	5.7	64	0.00
21	Hexachlorobenzene	0.236	0.220	6.8	62	0.00
22	Pentachlorophenol	0.037	0.048	-29.7#	93	-0.01
23	Phenanthrene	1.031	0.977	5.2	63	0.00
24	Anthracene	0.890	0.802	9.9	63	0.00
25 SURR	Fluoranthene-d10	5.063	4.841	4.4	65	0.00
26	Fluoranthene	1.313	1.245	5.2	65	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	60	0.00
28	Pvrene	1.016	0.987	2.9	65	0.00
29 S	Terphenyl-d14	0.817	0.785	3.9	65	0.00
30	Benzo(a)anthracene	1.075	1.039	3.3	65	-0.01
31	Chrysene	1.202	1.172	2.5	64	0.00
32	Bis(2-ethylhexyl)phthalate	0.787	0.788	-0.1	70	0.00
33	Indeno(1,2,3-cd)pyrene	1.044	0.905	13.3	57	0.00
34 I	Perylene-d12	1.000	1.000	0.0	59	0.00
35	Benzo(b)fluoranthene	1.098	1.020	7.1	62	0.00
36	Benzo(k)fluoranthene	1.217	1.174	3.5	64	0.00
37 C	Benzo(a)pyrene	0.994	0.932	6.2	62	0.00
38	Dibenzo(a,h)anthracene	0.979	0.857	12.5	58	-0.01
39	Benzo(a,h,i)perylene	1.024	0.907	11.4	57	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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