

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE102418\
 Data File : BE098047.D
 Acq On : 24 Oct 2018 20:26
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 25 04:03:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE102418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 24 13:49:35 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	88	0.00
2	1,4-Dioxane	0.616	0.611	0.8	93	0.00
3	n-Nitrosodimethylamine	0.821	0.753	8.3	85	0.00
4 S	2-Fluorophenol	1.211	1.115	7.9	84	0.00
5 S	Phenol-d6	2.017	1.896	6.0	85	0.00
6	bis(2-Chloroethyl)ether	1.413	1.173	17.0	75	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
8 S	Nitrobenzene-d5	0.425	0.394	7.3	86	0.00
9	Naphthalene	3.973	3.669	7.7	85	0.00
10	Hexachlorobutadiene	0.269	0.260	3.3	91	0.00
11 SURR	2-Methylnaphthalene-d10	0.710	0.652	8.2	87	0.00
12	2-Methylnaphthalene	0.719	0.664	7.6	84	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
14 S	2,4,6-Tribromophenol	0.201	0.180	10.4	88	0.00
15 S	2-Fluorobiphenyl	1.618	1.437	11.2	87	0.00
16	Acenaphthylene	1.811	1.679	7.3	89	-0.01
17	Acenaphthene	1.095	1.013	7.5	87	0.00
18	Fluorene	1.476	1.355	8.2	88	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
20	4-Bromophenyl-phenylether	0.241	0.218	9.5	87	0.00
21	Hexachlorobenzene	0.252	0.230	8.7	90	-0.01
22	Pentachlorophenol	0.021	0.020	4.8	98	0.00
23	Phenanthrene	0.982	0.880	10.4	88	-0.01
24	Anthracene	0.964	0.852	11.6	89	0.00
25 SURR	Fluoranthene-d10	5.129	4.564	11.0	88	0.00
26	Fluoranthene	1.224	1.080	11.8	87	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	90	0.00
28	Pvrene	1.105	0.976	11.7	86	0.00
29 S	Terphenyl-d14	0.943	0.832	11.8	88	0.00
30	Benzo(a)anthracene	1.182	1.077	8.9	89	-0.01
31	Chrysene	1.148	1.015	11.6	86	0.00
32	Bis(2-ethylhexyl)phthalate	2.807	2.393	14.7	83	0.00
33	Indeno(1,2,3-cd)pyrene	1.255	1.129	10.0	87	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	92	-0.01
35	Benzo(b)fluoranthene	1.111	0.999	10.1	87	-0.01
36	Benzo(k)fluoranthene	1.155	1.022	11.5	87	0.00
37 C	Benzo(a)pyrene	1.092	0.972	11.0	88	-0.01
38	Dibenzo(a,h)anthracene	1.064	0.947	11.0	86	0.00
39	Benzo(a,h,i)perylene	1.055	0.939	11.0	87	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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