

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE110718\
 Data File : BE098165.D
 Acq On : 7 Nov 2018 9:39
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Nov 07 10:15:13 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE110118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 01 13:00:52 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	89	0.00
2	1,4-Dioxane	0.596	0.628	-5.4	94	0.00
3	n-Nitrosodimethylamine	0.797	0.887	-11.3	106	0.00
4 S	2-Fluorophenol	1.187	1.236	-4.1	91	0.00
5 S	Phenol-d6	1.955	1.939	0.8	88	0.00
6	bis(2-Chloroethyl)ether	1.431	1.438	-0.5	93	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
8 S	Nitrobenzene-d5	0.429	0.419	2.3	90	-0.01
9	Naphthalene	3.998	4.076	-2.0	93	0.00
10	Hexachlorobutadiene	0.262	0.269	-2.7	94	0.00
11 SURR	2-Methylnaphthalene-d10	0.705	0.741	-5.1	94	0.00
12	2-Methylnaphthalene	0.714	0.729	-2.1	94	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	81	-0.01
14 S	2,4,6-Tribromophenol	0.189	0.159	15.9	81	-0.01
15 S	2-Fluorobiphenyl	1.645	1.706	-3.7	93	0.00
16	Acenaphthylene	1.825	1.922	-5.3	92	0.00
17	Acenaphthene	1.096	1.152	-5.1	89	0.00
18	Fluorene	1.454	1.509	-3.8	91	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
20	4-Bromophenyl-phenylether	0.244	0.255	-4.5	93	-0.01
21	Hexachlorobenzene	0.255	0.264	-3.5	93	0.00
22	Pentachlorophenol	0.057	0.053	7.0	91	0.00
23	Phenanthrene	0.991	1.014	-2.3	92	-0.01
24	Anthracene	0.942	0.960	-1.9	93	-0.01
25 SURR	Fluoranthene-d10	5.159	5.247	-1.7	93	0.00
26	Fluoranthene	1.241	1.274	-2.7	94	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	84	0.00
28	Pyrene	1.092	1.122	-2.7	94	0.00
29 S	Terphenyl-d14	0.930	0.942	-1.3	93	0.00
30	Benzo(a)anthracene	1.169	1.160	0.8	91	0.00
31	Chrysene	1.156	1.174	-1.6	92	0.00
32	Bis(2-ethylhexyl)phthalate	4.140	2.583	37.6#	66	0.00
33	Indeno(1,2,3-cd)pyrene	1.203	1.152	4.2	87	0.00
34 I	Perylene-d12	1.000	1.000	0.0	83	0.00
35	Benzo(b)fluoranthene	1.127	1.151	-2.1	90	0.00
36	Benzo(k)fluoranthene	1.166	1.212	-3.9	92	0.00
37 C	Benzo(a)pyrene	1.101	1.117	-1.5	89	0.00
38	Dibenzo(a,h)anthracene	1.027	0.984	4.2	84	0.00
39	Benzo(g,h,i)perylene	1.022	1.003	1.9	86	0.00

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