

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN073118\
 Data File : BN002201.D
 Acq On : 31 Jul 2018 08:32
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDC00.4

Quant Time: Jul 31 09:07:34 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 16:17:29 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	96	0.00
2	1,4-Dioxane	0.450	0.409	9.1	97	0.00
3	n-Nitrosodimethylamine	0.302	0.218	27.8#	78	0.00
4 S	2-Fluorophenol	0.933	0.847	9.2	92	0.00
5 S	Phenol-d6	1.135	1.024	9.8	94	0.00
6	bis(2-Chloroethyl)ether	1.091	0.974	10.7	87	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
8 S	Nitrobenzene-d5	0.301	0.285	5.3	96	0.00
9	Naphthalene	0.902	0.893	1.0	100	0.00
10	Hexachlorobutadiene	0.195	0.190	2.6	98	0.00
11 SURR	2-Methylnaphthalene-d10	0.589	0.571	3.1	99	0.00
12	2-Methylnaphthalene	0.610	0.599	1.8	98	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
14 S	2,4,6-Tribromophenol	0.171	0.143	16.4	82	0.00
15 S	2-Fluorobiphenyl	1.536	1.476	3.9	97	0.00
16	Acenaphthylene	1.709	1.598	6.5	94	0.00
17	Acenaphthene	1.099	1.071	2.5	97	0.00
18	Fluorene	1.352	1.290	4.6	92	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
20	4-Bromophenyl-phenylether	0.226	0.216	4.4	88	0.00
21	Hexachlorobenzene	0.250	0.250	0.0	90	0.00
22	Pentachlorophenol	0.046	0.038	17.4	74	0.00
23	Phenanthrene	0.995	0.952	4.3	84	0.00
24	Anthracene	0.823	0.749	9.0	81	0.00
25 SURR	Fluoranthene-d10	0.988	0.883	10.6	74	0.00
26	Fluoranthene	1.083	0.964	11.0	73	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	63	0.00
28	Pvrene	1.183	1.265	-6.9	72	0.00
29 S	Terphenyl-d14	0.855	0.888	-3.9	70	0.00
30	Benzo(a)anthracene	1.006	0.935	7.1	61	0.00
31	Chrysene	1.126	1.100	2.3	64	0.00
32	Bis(2-ethylhexyl)phthalate	0.460	0.394	14.3	55	-0.01
33	Indeno(1,2,3-cd)pyrene	1.061	1.088	-2.5	76	0.00
34 I	Perylene-d12	1.000	1.000	0.0	68	0.00
35	Benzo(b)fluoranthene	1.176	1.122	4.6	66	0.00
36	Benzo(k)fluoranthene	1.221	1.166	4.5	68	0.00
37 C	Benzo(a)pyrene	1.066	1.046	1.9	70	0.00
38	Dibenzo(a,h)anthracene	1.037	1.022	1.4	77	0.00
39	Benzo(a,h,i)perylene	1.060	1.048	1.1	76	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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