

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031120\
 Data File : BN010204.D
 Acq On : 11 Mar 2020 22:22
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 12 08:02:36 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN021920.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 11 17:37:44 2020
 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	62	0.02
2	1,4-Dioxane	0.410	0.486	-18.5	74	0.00
3	n-Nitrosodimethylamine	0.673	1.315	-95.4#	138	0.05
4 S	2-Fluorophenol	0.889	0.824	7.3	60	0.03
5 S	Phenol-d6	1.082	1.400	-29.4#	85	0.06
6	bis(2-Chloroethyl)ether	1.052	0.774	26.4#	48#	0.05
7 I	Naphthalene-d8	1.000	1.000	0.0	80	0.03
8 S	Nitrobenzene-d5	0.333	0.286	14.1	73	0.06
9	Naphthalene	1.091	1.070	1.9	81	0.03
10	Hexachlorobutadiene	0.240	0.249	-3.8	83	0.00
11 SURR	2-Methylnaphthalene-d10	0.774	0.621	19.8	66	0.03
12	2-Methylnaphthalene	0.748	0.578	22.7	64	0.04
13 I	Acenaphthene-d10	1.000	1.000	0.0	63	0.01
14 S	2,4,6-Tribromophenol	0.154	0.170	-10.4	78	0.02
15 S	2-Fluorobiphenyl	1.510	1.532	-1.5	67	0.02
16	Acenaphthylene	1.705	1.662	2.5	69	0.01
17	Acenaphthene	1.199	1.214	-1.3	67	0.01
18	Fluorene	1.582	1.576	0.4	65	0.02
19 I	Phenanthrene-d10	1.000	1.000	0.0	61	0.02
20	4-Bromophenyl-phenylether	0.279	0.159	43.0#	37#	0.00
21	Hexachlorobenzene	0.249	0.259	-4.0	66	0.01
22	Pentachlorophenol	0.064	0.056	12.5	72	0.02
23	Phenanthrene	1.190	1.093	8.2	61	0.01
24	Anthracene	1.006	0.925	8.1	62	0.02
25 SURR	Fluoranthene-d10	1.384	1.365	1.4	65	0.01
26	Fluoranthene	1.425	1.395	2.1	64	0.01
27 I	Chrysene-d12	1.000	1.000	0.0	62	0.01
28	Pvrene	1.519	1.550	-2.0	65	0.01
29 S	Terphenyl-d14	0.956	1.034	-8.2	69	0.02
30	Benzo(a)anthracene	1.327	1.106	16.7	58	0.01
31	Chrysene	1.536	1.706	-11.1	71	0.01
32	Bis(2-ethylhexyl)phthalate	0.648	0.628	3.1	67	0.00
33	Indeno(1,2,3-cd)pyrene	1.599	1.735	-8.5	67	0.01
34 I	Perylene-d12	1.000	1.000	0.0	67	0.00
35	Benzo(b)fluoranthene	1.462	1.303	10.9	65	0.01
36	Benzo(k)fluoranthene	1.592	1.616	-1.5	72	0.01
37 C	Benzo(a)pyrene	1.334	1.345	-0.8	70	0.01
38	Dibenzo(a,h)anthracene	1.314	1.210	7.9	66	0.03
39	Benzo(a,h,i)perylene	1.400	1.483	-5.9	75	0.02

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