

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032319\
 Data File : BN004933.D
 Acq On : 23 Mar 2019 22:36
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDC00.4

Quant Time: Mar 24 02:13:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN032019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 20 16:06:20 2019
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	90	0.00
2	1,4-Dioxane	0.400	0.440	-10.0	99	-0.02
3	n-Nitrosodimethylamine	0.400	0.453	-13.2	111	0.00
4 S	2-Fluorophenol	0.400	0.492	-23.0	109	0.00
5 S	Phenol-d6	0.400	0.504	-26.0#	115	0.00
6	bis(2-Chloroethyl)ether	0.400	0.460	-15.0	107	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	90	-0.01
8 S	Nitrobenzene-d5	0.400	0.485	-21.2	110	-0.03
9	Naphthalene	0.400	0.435	-8.7	97	-0.03
10	Hexachlorobutadiene	0.400	0.410	-2.5	91	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.396	1.0	89	-0.01
12	2-Methylnaphthalene	0.400	0.402	-0.5	90	-0.01
13 I	Acenaphthene-d10	0.400	0.400	0.0	77	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.431	-7.7	89	-0.01
15 S	2-Fluorobiphenyl	0.400	0.457	-14.2	87	-0.01
16	Acenaphthylene	0.400	0.428	-7.0	87	-0.01
17	Acenaphthene	0.400	0.423	-5.7	81	-0.01
18	Fluorene	0.400	0.395	1.3	75	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	70	-0.01
20	4-Bromophenyl-phenylether	0.400	0.406	-1.5	71	-0.01
21	Hexachlorobenzene	0.400	0.393	1.8	68	-0.01
22	Pentachlorophenol	0.400	0.613	-53.2#	123	-0.01
23	Phenanthrene	0.400	0.418	-4.5	72	-0.01
24	Anthracene	0.400	0.421	-5.2	75	-0.01
25 SURR	Fluoranthene-d10	0.400	0.425	-6.2	75	-0.01
26	Fluoranthene	0.400	0.421	-5.2	73	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	76	-0.01
28	Pvrene	0.400	0.396	1.0	74	-0.02
29 S	Terphenyl-d14	0.400	0.407	-1.7	76	-0.01
30	Benzo(a)anthracene	0.400	0.423	-5.7	81	-0.02
31	Chrysene	0.400	0.414	-3.5	76	-0.02
32	Bis(2-ethylhexyl)phthalate	0.400	0.438	-9.5	90	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.530	-32.5#	97	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	94	-0.02
35	Benzo(b)fluoranthene	0.400	0.373	6.8	86	-0.01
36	Benzo(k)fluoranthene	0.400	0.364	9.0	87	-0.01
37 C	Benzo(a)pyrene	0.400	0.398	0.5	94	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.413	-3.2	96	-0.02
39	Benzo(a,h,i)perylene	0.400	0.419	-4.7	97	-0.03

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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