

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN032419\
 Data File : BN004952.D
 Acq On : 24 Mar 2019 22:53
 Operator : JU/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 25 04:13:05 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	261	-0.02
2	1,4-Dioxane	0.400	0.426	-6.5	275	-0.02
3	n-Nitrosodimethylamine	0.400	0.488	-22.0	344	-0.02
4 S	2-Fluorophenol	0.400	0.484	-21.0	311	0.00
5 S	Phenol-d6	0.400	0.490	-22.5	322	0.00
6	bis(2-Chloroethyl)ether	0.400	0.450	-12.5	302	-0.02
7 I	Naphthalene-d8	0.400	0.400	0.0	254	-0.03
8 S	Nitrobenzene-d5	0.400	0.499	-24.7	320	-0.03
9	Naphthalene	0.400	0.436	-9.0	273	-0.03
10	Hexachlorobutadiene	0.400	0.412	-3.0	259	-0.03
11 SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	253	-0.02
12	2-Methylnaphthalene	0.400	0.403	-0.8	254	-0.02
13 I	Acenaphthene-d10	0.400	0.400	0.0	206	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.398	0.5	218	-0.01
15 S	2-Fluorobiphenyl	0.400	0.473	-18.2	239	-0.02
16	Acenaphthylene	0.400	0.488	-22.0	264	-0.02
17	Acenaphthene	0.400	0.440	-10.0	225	-0.01
18	Fluorene	0.400	0.405	-1.3	204	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	166	-0.01
20	4-Bromophenyl-phenylether	0.400	0.458	-14.5	191	-0.01
21	Hexachlorobenzene	0.400	0.413	-3.2	169	-0.01
22	Pentachlorophenol	0.400	0.534	-33.5#	256	-0.01
23	Phenanthrene	0.400	0.415	-3.7	171	-0.01
24	Anthracene	0.400	0.444	-11.0	188	-0.01
25 SURR	Fluoranthene-d10	0.400	0.369	7.8	154	-0.02
26	Fluoranthene	0.400	0.363	9.3	150	-0.02
27 I	Chrysene-d12	0.400	0.400	0.0	123	-0.01
28	Pyrene	0.400	0.481	-20.2	146	-0.02
29 S	Terphenyl-d14	0.400	0.509	-27.2#	155	-0.02
30	Benzo(a)anthracene	0.400	0.452	-13.0	141	-0.02
31	Chrysene	0.400	0.406	-1.5	121	-0.02
32	Bis(2-ethylhexyl)phthalate	0.400	0.785	-96.3#	263	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.371	7.3	111	-0.02
34 I	Perylene-d12	0.400	0.400	0.0	118	-0.02
35	Benzo(b)fluoranthene	0.400	0.418	-4.5	121	-0.01
36	Benzo(k)fluoranthene	0.400	0.394	1.5	117	-0.01
37 C	Benzo(a)pyrene	0.400	0.404	-1.0	119	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.376	6.0	109	-0.02
39	Benzo(g,h,i)perylene	0.400	0.377	5.8	109	-0.03

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

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