

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070919\
 Data File : BN006768.D
 Acq On : 09 Jul 2019 15:58
 Operator : HP/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 ICVBN070919

Quant Time: Jul 09 17:36:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN070919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 09 17:31:09 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	98	0.00
2	1,4-Dioxane	0.657	0.608	7.5	92	0.00
3	n-Nitrosodimethylamine	0.510	0.424	16.9	94	0.00
4 S	2-Fluorophenol	0.889	0.774	12.9	87	0.00
5 S	Phenol-d6	1.223	1.022	16.4	86	0.00
6	bis(2-Chloroethyl)ether	1.361	1.130	17.0	87	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	99	0.01
8 S	Nitrobenzene-d5	0.273	0.234	14.3	87	0.01
9	Naphthalene	1.071	1.028	4.0	88	0.00
10	Hexachlorobutadiene	0.232	0.226	2.6	87	0.01
11 SURR	2-Methylnaphthalene-d10	0.598	0.552	7.7	86	0.00
12	2-Methylnaphthalene	0.685	0.652	4.8	89	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
14 S	2,4,6-Tribromophenol	0.187	0.155	17.1	84	0.01
15 S	2-Fluorobiphenyl	1.490	1.416	5.0	88	0.00
16	Acenaphthylene	1.889	1.741	7.8	88	0.00
17	Acenaphthene	1.268	1.217	4.0	87	0.01
18	Fluorene	1.571	1.462	6.9	85	0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.01
20	4-Bromophenyl-phenylether	0.226	0.207	8.4	84	0.01
21	Hexachlorobenzene	0.295	0.284	3.7	85	0.01
22	Pentachlorophenol	0.039	0.035	10.3	93	0.00
23	Phenanthrene	1.109	1.037	6.5	85	0.01
24	Anthracene	0.961	0.850	11.6	84	0.00
25 SURR	Fluoranthene-d10	1.208	1.109	8.2	83	0.00
26	Fluoranthene	1.392	1.316	5.5	85	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	93	0.02
28	Pvrene	1.390	1.376	1.0	84	0.00
29 S	Terphenyl-d14	0.881	0.865	1.8	85	0.00
30	Benzo(a)anthracene	1.233	1.115	9.6	81	0.02
31	Chrysene	1.382	1.395	-0.9	84	0.02
32	Bis(2-ethylhexyl)phthalate	0.574	0.477	16.9	81	0.02
33	Indeno(1,2,3-cd)pyrene	1.691	1.707	-0.9	82	0.05
34 I	Perylene-d12	1.000	1.000	0.0	92	0.04
35	Benzo(b)fluoranthene	1.257	1.206	4.1	86	0.03
36	Benzo(k)fluoranthene	1.411	1.432	-1.5	87	0.03
37 C	Benzo(a)pyrene	1.219	1.205	1.1	85	0.04
38	Dibenzo(a,h)anthracene	1.229	1.189	3.3	83	0.05
39	Benzo(a,h,i)perylene	1.288	1.261	2.1	82	0.05

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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