

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090820\
 Data File : BN011729.D
 Acq On : 08 Sep 2020 10:26
 Operator : CG/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTDCCC0.4

Quant Time: Sep 08 14:38:02 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN090820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 08 14:34:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	1757	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	6345	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	3362	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	7212	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	5967	0.40	ng/ul	0.00
20) Perylene-d12	23.57	264	5783	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	4215	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	8151	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.57	128	8102	0.431	ng/ul	99
5) 2-Methylnaphthalene	12.19	142	5434	0.434	ng/ul	100
7) Acenaphthylene	14.09	152	7127	0.440	ng/ul	100
8) Acenaphthene	14.44	153	5679	0.436	ng/ul	99
9) Fluorene	15.43	166	6501	0.432	ng/ul	100
11) Pentachlorophenol	16.77	266	925	0.425	ng/ul	98
12) Phenanthrene	17.16	178	10619	0.438	ng/ul	99
13) Anthracene	17.25	178	9422	0.436	ng/ul	99
15) Fluoranthene	19.18	202	11712	0.434	ng/ul	100
17) Pyrene	19.54	202	11622	0.439	ng/ul	98
18) Benzo(a)anthracene	21.30	228	10079	0.434	ng/ul	100
19) Chrysene	21.35	228	11410	0.440	ng/ul	98
21) Benzo(b)fluoranthene	22.89	252	11388	0.420	ng/ul	99
22) Benzo(k)fluoranthene	22.94	252	11779	0.424	ng/ul	100
23) Benzo(a)pyrene	23.47	252	9791	0.418	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.83	276	13522	0.440	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.84	278	10918	0.435	ng/ul	99
26) Benzo(g,h,i)perylene	26.52	276	11348	0.422	ng/ul	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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