

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120120\
 Data File : BN012770.D
 Acq On : 30 Nov 2020 18:07
 Operator : CG/JU
 Sample : PB133258BL
 Misc : MDL-WATER-BL-07
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK258

Quant Time: Dec 01 00:53:55 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN112520.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 30 17:13:00 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	129001	20.00	ng/ul	0.00
20) Naphthalene-d8	10.19	136	530189	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.08	164	349651	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.83	188	740096	20.00	ng/ul	0.00
79) Chrysene-d12	21.06	240	741504	20.00	ng/ul	0.00
88) Perylene-d12	23.17	264	860365	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.96	96	22881	6.39	ng/uL	0.00
4) Pyridine-d5	3.36	84	338359	34.95	ng/ul	0.00
7) Phenol-d5	6.60	99	392727	34.01	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.77	67	251296	34.89	ng/ul	0.00
11) 2-Chlorophenol-d4	6.95	132	321752	35.44	ng/ul	0.00
15) 4-Methylphenol-d8	8.13	113	321949	34.23	ng/ul	0.00
21) Nitrobenzene-d5	8.57	128	157096	37.03	ng/ul	0.00
24) 2-Nitrophenol-d4	9.28	143	175245	37.42	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.81	165	311605	35.03	ng/ul	0.00
31) 4-Chloroaniline-d4	10.33	131	430212	34.57	ng/ul	0.00
46) Dimethylphthalate-d6	13.50	166	1014661	35.69	ng/ul	0.00
49) Acenaphthylene-d8	13.77	160	1134305	34.19	ng/ul	0.00
54) 4-Nitrophenol-d4	14.30	143	157992	31.41	ng/ul	0.00
60) Fluorene-d10	15.08	176	865487	36.06	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.22	200	143421	28.34	ng/ul	0.00
73) Anthracene-d10	16.93	188	1331967	36.01	ng/ul	0.00
81) Pyrene-d10	19.25	212	1506027	40.74	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.04	264	1682233	36.19	ng/ul	0.00

Target Compounds Ovalue

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

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