

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN120120\
 Data File : BN012768.D
 Acq On : 30 Nov 2020 16:54
 Operator : CG/JU
 Sample : PB133258BL
 Misc : MDL-WATER-BL-06
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK258

Quant Time: Nov 30 17:28:28 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	157666	20.00	ng/ul	0.00
20) Naphthalene-d8	10.19	136	645423	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.08	164	425084	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.83	188	895499	20.00	ng/ul	0.00
79) Chrysene-d12	21.06	240	960739	20.00	ng/ul	0.00
88) Perylene-d12	23.17	264	1062417	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.96	96	28482	6.50	ng/uL	0.00
4) Pyridine-d5	3.36	84	402906	34.05	ng/ul	0.00
7) Phenol-d5	6.61	99	474990	33.65	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.77	67	307868	34.97	ng/ul	0.00
11) 2-Chlorophenol-d4	6.96	132	384679	34.67	ng/ul	0.00
15) 4-Methylphenol-d8	8.13	113	385805	33.56	ng/ul	0.00
21) Nitrobenzene-d5	8.57	128	187133	36.23	ng/ul	0.00
24) 2-Nitrophenol-d4	9.29	143	217491	38.15	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.81	165	374645	34.59	ng/ul	0.00
31) 4-Chloroaniline-d4	10.33	131	513106	33.87	ng/ul	0.00
46) Dimethylphthalate-d6	13.50	166	1301777	37.66	ng/ul	0.00
49) Acenaphthylene-d8	13.77	160	1413506	35.05	ng/ul	0.00
54) 4-Nitrophenol-d4	14.30	143	201472	32.95	ng/ul	0.00
60) Fluorene-d10	15.08	176	1083522	37.13	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.22	200	182160	29.75	ng/ul	0.00
73) Anthracene-d10	16.93	188	1662796	37.16	ng/ul	0.00
81) Pyrene-d10	19.24	212	1844413	38.51	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.04	264	2199205	38.31	ng/ul	0.00

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

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