

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN112520\
 Data File : BN012738.D
 Acq On : 25 Nov 2020 19:06
 Operator : CG/JU
 Sample : PB133233BL
 Misc : MDL-WATER-BL-01
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK233

Quant Time: Nov 27 11:01:34 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN112520.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 25 18:19:04 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.42	152	119478	20.00	ng/ul	0.00
20) Naphthalene-d8	10.19	136	487639	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.08	164	312516	20.00	ng/ul	0.00
64) Phenanthrene-d10	16.84	188	681755	20.00	ng/ul	0.00
79) Chrysene-d12	21.06	240	665549	20.00	ng/ul	0.00
88) Perylene-d12	23.17	264	778875	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.96	96	23357	7.04	ng/uL	0.00
4) Pyridine-d5	3.36	84	331674	36.99	ng/ul	0.00
7) Phenol-d5	6.61	99	396604	37.08	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	6.78	67	249490	37.40	ng/ul	0.00
11) 2-Chlorophenol-d4	6.96	132	319270	37.97	ng/ul	0.00
15) 4-Methylphenol-d8	8.13	113	321573	36.91	ng/ul	0.00
21) Nitrobenzene-d5	8.58	128	154822	39.68	ng/ul	0.00
24) 2-Nitrophenol-d4	9.29	143	179319	41.63	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.82	165	296718	36.26	ng/ul	0.00
31) 4-Chloroaniline-d4	10.34	131	426807	37.29	ng/ul	0.00
46) Dimethylphthalate-d6	13.50	166	1036090	40.77	ng/ul	0.00
49) Acenaphthylene-d8	13.77	160	1138541	38.40	ng/ul	0.00
54) 4-Nitrophenol-d4	14.30	143	155059	34.49	ng/ul	0.00
60) Fluorene-d10	15.09	176	853517	39.78	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.22	200	153719	32.98	ng/ul	0.00
73) Anthracene-d10	16.93	188	1312373	38.52	ng/ul	0.00
81) Pyrene-d10	19.25	212	1477714	44.54	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.04	264	1653566	39.30	ng/ul	0.00

Target Compounds Ovalue

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

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