

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG113020\
 Data File : BG047483.D
 Acq On : 30 Nov 2020 10:46
 Operator : CG/JU
 Sample : PB133258BL
 Misc : MDL-WATER-BL-06
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK258

Quant Time: Nov 30 11:24:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG111820.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 30 10:51:47 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	33298	20.00	ng/ul	0.00
20) Naphthalene-d8	11.09	136	127248	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.88	164	101047	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	259070	20.00	ng/ul	0.00
79) Chrysene-d12	21.90	240	278181	20.00	ng/ul	0.00
88) Perylene-d12	25.31	264	277371	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.61	96	4640	5.52	ng/uL	0.00
4) Pyridine-d5	4.03	84	76088	29.49	ng/ul	0.00
7) Phenol-d5	7.40	99	99458	31.17	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.57	67	54654	29.27	ng/ul	0.00
11) 2-Chlorophenol-d4	7.79	132	73183	33.61	ng/ul	0.00
15) 4-Methylphenol-d8	8.96	113	75325	30.40	ng/ul	0.00
21) Nitrobenzene-d5	9.44	128	37444	35.45	ng/ul	0.00
24) 2-Nitrophenol-d4	10.17	143	43231	39.11	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.71	165	84800	33.35	ng/ul	0.00
31) 4-Chloroaniline-d4	11.22	131	104031	33.12	ng/ul	0.00
46) Dimethylphthalate-d6	14.27	166	308957	36.56	ng/ul	0.00
49) Acenaphthylene-d8	14.58	160	325016	33.71	ng/ul	0.00
54) 4-Nitrophenol-d4	15.04	143	39743	32.21	ng/ul	0.00
60) Fluorene-d10	15.86	176	262582	34.39	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.96	200	46451	33.84	ng/ul	0.00
73) Anthracene-d10	17.72	188	455152	36.29	ng/ul	0.00
81) Pyrene-d10	19.98	212	535748	34.13	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.07	264	555897	36.35	ng/ul	0.01

Target Compounds Ovalue

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

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