

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG112520\
 Data File : BG047469.D
 Acq On : 25 Nov 2020 15:39
 Operator : CG/JU
 Sample : PB133233BL
 Misc : MDL-WATER-BL02
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK233

Manual Integrations
 APPROVED

mohammad
 11/27/2020 11:45:41 AM

Quant Time: Nov 25 16:24:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG111820.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 25 12:54:28 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	27172	20.00	ng/ul	0.00
20) Naphthalene-d8	11.09	136	106496	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.88	164	84536	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.61	188	208735	20.00	ng/ul	0.00
79) Chrysene-d12	21.90	240	215299	20.00	ng/ul	0.00
88) Perylene-d12	25.29	264	219079	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.61	96	5009m	7.30	ng/uL	0.00
4) Pyridine-d5	4.04	84	80236	38.11	ng/ul	0.00
7) Phenol-d5	7.40	99	108956	41.85	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.58	67	58213	38.20	ng/ul	0.00
11) 2-Chlorophenol-d4	7.80	132	78695	44.29	ng/ul	0.00
15) 4-Methylphenol-d8	8.96	113	79460	39.30	ng/ul	0.00
21) Nitrobenzene-d5	9.44	128	39272	44.42	ng/ul	0.00
24) 2-Nitrophenol-d4	10.16	143	48120	52.02	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.71	165	93214	43.80	ng/ul	0.00
31) 4-Chloroaniline-d4	11.22	131	110991	42.23	ng/ul	0.00
46) Dimethylphthalate-d6	14.27	166	323689	45.79	ng/ul	0.00
49) Acenaphthylene-d8	14.58	160	347393	43.07	ng/ul	0.00
54) 4-Nitrophenol-d4	15.04	143	45819	44.38	ng/ul	0.00
60) Fluorene-d10	15.86	176	278917	43.66	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.96	200	49657	44.90	ng/ul	0.00
73) Anthracene-d10	17.71	188	471946	46.70	ng/ul	0.00
81) Pyrene-d10	19.98	212	556903	45.84	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.06	264	552196	45.72	ng/ul	0.00

Target Compounds Ovalue

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

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