

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120820\
 Data File : BM028081.D
 Acq On : 08 Dec 2020 21:11
 Operator : JU/CG
 Sample : PB133258BL
 Misc : MDL-WATER-BL-01
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK258

Quant Time: Dec 09 03:54:30 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120820.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 09 03:34:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	153664	20.00	ng/ul	0.00
20) Naphthalene-d8	11.12	136	624159	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.90	164	364399	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	753127	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	736281	20.00	ng/ul	-0.01
88) Perylene-d12	24.13	264	733276	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	29279	7.28	ng/uL	0.00
4) Pyridine-d5	3.92	84	424249	38.60	ng/ul	0.00
7) Phenol-d5	7.41	99	509308	38.83	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.59	67	323099	38.77	ng/ul	0.00
11) 2-Chlorophenol-d4	7.80	132	434008	39.74	ng/ul	0.00
15) 4-Methylphenol-d8	8.98	113	404119	37.75	ng/ul	0.00
21) Nitrobenzene-d5	9.46	128	200696	39.85	ng/ul	0.00
24) 2-Nitrophenol-d4	10.19	143	210966	40.68	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.72	165	388373	38.11	ng/ul	0.00
31) 4-Chloroaniline-d4	11.24	131	570521	40.13	ng/ul	0.00
46) Dimethylphthalate-d6	14.30	166	1215169	41.94	ng/ul	0.00
49) Acenaphthylene-d8	14.60	160	1384714	39.33	ng/ul	0.00
54) 4-Nitrophenol-d4	15.07	143	205956	37.73	ng/ul	0.00
60) Fluorene-d10	15.89	176	1000535	39.70	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.99	200	153667	33.40	ng/ul	0.00
73) Anthracene-d10	17.72	188	1541910	41.13	ng/ul	0.00
81) Pyrene-d10	19.97	212	1665423	41.66	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.98	264	1696635	42.26	ng/ul	0.00

Target Compounds Ovalue

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

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