

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120720\
 Data File : BM028059.D
 Acq On : 07 Dec 2020 15:19
 Operator : JU/CG
 Sample : PB133234BL
 Misc : MDL-WATER-BL-02
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK234

Quant Time: Dec 07 16:04:39 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120320.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 07 13:18:59 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	232375	20.00	ng/ul	0.00
20) Naphthalene-d8	11.12	136	948373	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.91	164	539684	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	1052118	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	925647	20.00	ng/ul	0.00
88) Perylene-d12	24.14	264	899083	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	35667	5.73	ng/uL	0.00
4) Pyridine-d5	3.93	84	495555	29.52	ng/ul	0.00
7) Phenol-d5	7.42	99	601418	31.09	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.59	67	382841	31.84	ng/ul	0.00
11) 2-Chlorophenol-d4	7.80	132	505695	31.38	ng/ul	0.00
15) 4-Methylphenol-d8	8.99	113	477413	31.29	ng/ul	0.00
21) Nitrobenzene-d5	9.46	128	237811	32.30	ng/ul	0.00
24) 2-Nitrophenol-d4	10.19	143	249188	33.87	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.73	165	456543	30.60	ng/ul	0.00
31) 4-Chloroaniline-d4	11.25	131	662338	32.45	ng/ul	0.00
46) Dimethylphthalate-d6	14.30	166	1357529	34.00	ng/ul	0.00
49) Acenaphthylene-d8	14.60	160	1573527	30.70	ng/ul	0.00
54) 4-Nitrophenol-d4	15.07	143	224059	29.10	ng/ul	0.00
60) Fluorene-d10	15.89	176	1125609	31.45	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.99	200	167982	27.51	ng/ul	0.00
73) Anthracene-d10	17.72	188	1673441	32.44	ng/ul	0.00
81) Pyrene-d10	19.97	212	1739475	35.80	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.99	264	1653916	34.55	ng/ul	0.00

Target Compounds Ovalue

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

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