

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120820\
 Data File : BM028083.D
 Acq On : 08 Dec 2020 22:23
 Operator : JU/CG
 Sample : PB133258BL
 Misc : MDL-WATER-BL-02
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK258

Quant Time: Dec 09 03:59:00 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	145575	20.00	ng/ul	0.00
20) Naphthalene-d8	11.12	136	586122	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.90	164	334716	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	680131	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	683282	20.00	ng/ul	-0.01
88) Perylene-d12	24.13	264	698365	20.00	ng/ul	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.46	96	29948	7.86	ng/uL	0.00
4) Pyridine-d5	3.93	84	405836	38.98	ng/ul	0.00
7) Phenol-d5	7.41	99	482610	38.84	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.59	67	302704	38.34	ng/ul	0.00
11) 2-Chlorophenol-d4	7.80	132	409030	39.54	ng/ul	0.00
15) 4-Methylphenol-d8	8.98	113	383705	37.83	ng/ul	0.00
21) Nitrobenzene-d5	9.46	128	188698	39.90	ng/ul	0.00
24) 2-Nitrophenol-d4	10.19	143	197633	40.58	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.72	165	362589	37.89	ng/ul	0.00
31) 4-Chloroaniline-d4	11.24	131	533564	39.97	ng/ul	0.00
46) Dimethylphthalate-d6	14.30	166	1100899	41.37	ng/ul	0.00
49) Acenaphthylene-d8	14.60	160	1264818	39.11	ng/ul	0.00
54) 4-Nitrophenol-d4	15.06	143	187569	37.41	ng/ul	0.00
60) Fluorene-d10	15.88	176	919692	39.73	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.99	200	137198	33.02	ng/ul	0.00
73) Anthracene-d10	17.72	188	1399090	41.33	ng/ul	0.00
81) Pyrene-d10	19.97	212	1530092	41.24	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.97	264	1613603	42.20	ng/ul	-0.01

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

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