

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120820\
 Data File : BP004196.D
 Acq On : 08 Dec 2020 12:57
 Operator : CG/JU
 Sample : PB133258BL
 Misc : MDL-WATER-BL-06
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK258

Quant Time: Dec 08 13:48:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP120420.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 08 06:09:39 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	388304	20.00	ng/ul	0.00
20) Naphthalene-d8	10.64	136	1624652	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.49	164	990147	20.00	ng/ul	-0.01
64) Phenanthrene-d10	17.25	188	2119650	20.00	ng/ul	0.00
79) Chrysene-d12	21.34	240	2082625	20.00	ng/ul	-0.01
88) Perylene-d12	23.70	264	2209947	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.34	96	78152	7.73	ng/uL	0.00
4) Pyridine-d5	3.74	84	1133978	39.49	ng/ul	0.00
7) Phenol-d5	7.00	99	1295557	37.32	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	7.17	67	822878	41.01	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.37	132	997655	36.93	ng/ul	-0.01
15) 4-Methylphenol-d8	8.54	113	997304	36.47	ng/ul	-0.01
21) Nitrobenzene-d5	9.00	128	536741	40.52	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	463630	38.34	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.25	165	948293	35.63	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.77	131	1457110	37.31	ng/ul	-0.01
46) Dimethylphthalate-d6	13.90	166	3242322	41.23	ng/ul	-0.01
49) Acenaphthylene-d8	14.18	160	3796873	39.29	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.67	143	438449	34.06	ng/ul	0.00
60) Fluorene-d10	15.49	176	2751623	40.01	ng/ul	-0.01
65) 4,6-Dinitro-2-methylphenol	15.60	200	284430	33.02	ng/ul	-0.01
73) Anthracene-d10	17.34	188	4288689	41.17	ng/ul	-0.01
81) Pyrene-d10	19.59	212	4796064	43.77	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.56	264	4983472	41.98	ng/ul	-0.01

Target Compounds Ovalue

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

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