

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120820\
 Data File : BP004198.D
 Acq On : 08 Dec 2020 14:05
 Operator : CG/JU
 Sample : PB133258BL
 Misc : MDL-WATER-BL-07
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK258

Quant Time: Dec 08 14:47:52 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	392957	20.00	ng/ul	0.00
20) Naphthalene-d8	10.64	136	1641895	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.49	164	997554	20.00	ng/ul	-0.01
64) Phenanthrene-d10	17.25	188	2149578	20.00	ng/ul	0.00
79) Chrysene-d12	21.34	240	2099755	20.00	ng/ul	0.00
88) Perylene-d12	23.71	264	2245050	20.00	ng/ul	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.34	96	81024	7.92	ng/uL	0.00
4) Pyridine-d5	3.73	84	1141970	39.30	ng/ul	-0.01
7) Phenol-d5	7.00	99	1307410	37.22	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	7.17	67	835310	41.13	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.37	132	1004988	36.76	ng/ul	-0.01
15) 4-Methylphenol-d8	8.54	113	1002399	36.22	ng/ul	-0.01
21) Nitrobenzene-d5	9.00	128	546400	40.82	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	464913	38.04	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.25	165	952084	35.40	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.77	131	1476919	37.42	ng/ul	-0.01
46) Dimethylphthalate-d6	13.90	166	3271681	41.30	ng/ul	-0.01
49) Acenaphthylene-d8	14.18	160	3840201	39.45	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.67	143	447746	34.53	ng/ul	-0.01
60) Fluorene-d10	15.49	176	2762824	39.88	ng/ul	-0.01
65) 4,6-Dinitro-2-methylphenol	15.60	200	288807	33.07	ng/ul	-0.01
73) Anthracene-d10	17.35	188	4325078	40.94	ng/ul	0.00
81) Pyrene-d10	19.59	212	4848773	43.89	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.56	264	5067173	42.02	ng/ul	-0.01

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

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