

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120520\
 Data File : BP004161.D
 Acq On : 05 Dec 2020 14:47
 Operator : CG/JU
 Sample : PB133233BL
 Misc : MDL-WATER-BL-01
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK233

Quant Time: Dec 07 02:46:03 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP120420.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 07 01:39:51 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	340517	20.00	ng/ul	0.00
20) Naphthalene-d8	10.65	136	1406178	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.49	164	889495	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.25	188	1983284	20.00	ng/ul	0.00
79) Chrysene-d12	21.35	240	2059014	20.00	ng/ul	0.00
88) Perylene-d12	23.71	264	2215209	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.35	96	55575	6.27	ng/uL	0.00
4) Pyridine-d5	3.74	84	780394	30.99	ng/ul	0.00
7) Phenol-d5	7.00	99	915688	30.08	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.18	67	561765	31.92	ng/ul	0.00
11) 2-Chlorophenol-d4	7.38	132	703014	29.68	ng/ul	0.00
15) 4-Methylphenol-d8	8.55	113	690184	28.78	ng/ul	0.00
21) Nitrobenzene-d5	9.00	128	350941	30.61	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	307465	29.37	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.26	165	645287	28.01	ng/ul	0.00
31) 4-Chloroaniline-d4	10.77	131	1005751	29.75	ng/ul	0.00
46) Dimethylphthalate-d6	13.90	166	2336812	33.08	ng/ul	0.00
49) Acenaphthylene-d8	14.19	160	2640081	30.41	ng/ul	0.00
54) 4-Nitrophenol-d4	14.67	143	320543	27.72	ng/ul	0.00
60) Fluorene-d10	15.49	176	1945548	31.49	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.60	200	197901	24.56	ng/ul	0.00
73) Anthracene-d10	17.35	188	3161861	32.44	ng/ul	0.00
81) Pyrene-d10	19.59	212	3650637	33.70	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.56	264	4008329	33.68	ng/ul	0.00

Target Compounds Ovalue

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

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