

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120720\
 Data File : BP004181.D
 Acq On : 07 Dec 2020 14:12
 Operator : CG/JU
 Sample : PB133234BL
 Misc : MDL-WATER-BL-04
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK234

Quant Time: Dec 07 15:33:52 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP120420.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Dec 07 12:11:35 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	419210	20.00	ng/ul	0.00
20) Naphthalene-d8	10.64	136	1769325	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.49	164	1086218	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.25	188	2392360	20.00	ng/ul	0.00
79) Chrysene-d12	21.34	240	2398273	20.00	ng/ul	-0.01
88) Perylene-d12	23.70	264	2595317	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.34	96	67713	6.21	ng/uL	0.00
4) Pyridine-d5	3.74	84	943726	30.44	ng/ul	0.00
7) Phenol-d5	7.00	99	1090127	29.09	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	7.18	67	701981	32.40	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.38	132	839437	28.78	ng/ul	-0.01
15) 4-Methylphenol-d8	8.54	113	831858	28.17	ng/ul	-0.01
21) Nitrobenzene-d5	9.00	128	443256	30.73	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	357929	27.18	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.26	165	783881	27.05	ng/ul	0.00
31) 4-Chloroaniline-d4	10.77	131	1223424	28.76	ng/ul	-0.01
46) Dimethylphthalate-d6	13.90	166	2774223	32.16	ng/ul	-0.01
49) Acenaphthylene-d8	14.18	160	3209443	30.28	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.67	143	361405	25.60	ng/ul	0.00
60) Fluorene-d10	15.49	176	2360185	31.28	ng/ul	-0.01
65) 4,6-Dinitro-2-methylphenol	15.60	200	227138	23.37	ng/ul	-0.01
73) Anthracene-d10	17.35	188	3749557	31.89	ng/ul	0.00
81) Pyrene-d10	19.59	212	4308538	34.15	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.56	264	4596941	32.97	ng/ul	-0.01

Target Compounds Ovalue

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

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