

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120720\
 Data File : BP004183.D
 Acq On : 07 Dec 2020 15:20
 Operator : CG/JU
 Sample : PB133234BL
 Misc : MDL-WATER-BL-05
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK234

Quant Time: Dec 07 16:03:04 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	456845	20.00	ng/ul	0.00
20) Naphthalene-d8	10.64	136	1908986	20.00	ng/ul	-0.01
38) Acenaphthene-d10	14.49	164	1174533	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.25	188	2567038	20.00	ng/ul	0.00
79) Chrysene-d12	21.34	240	2546782	20.00	ng/ul	-0.01
88) Perylene-d12	23.70	264	2753688	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.34	96	72964	6.14	ng/uL	0.00
4) Pyridine-d5	3.74	84	1019988	30.19	ng/ul	0.00
7) Phenol-d5	7.00	99	1191576	29.18	ng/ul	-0.01
9) Bis-(2-Chloroethyl)ether-d	7.18	67	763504	32.34	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.38	132	920231	28.96	ng/ul	-0.01
15) 4-Methylphenol-d8	8.54	113	913162	28.38	ng/ul	-0.01
21) Nitrobenzene-d5	9.00	128	486512	31.26	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	405176	28.51	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.26	165	852031	27.25	ng/ul	0.00
31) 4-Chloroaniline-d4	10.77	131	1320272	28.77	ng/ul	-0.01
46) Dimethylphthalate-d6	13.90	166	3015401	32.33	ng/ul	-0.01
49) Acenaphthylene-d8	14.18	160	3464343	30.22	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.67	143	398044	26.07	ng/ul	-0.01
60) Fluorene-d10	15.49	176	2539555	31.13	ng/ul	-0.01
65) 4,6-Dinitro-2-methylphenol	15.60	200	255754	24.52	ng/ul	-0.01
73) Anthracene-d10	17.35	188	4005229	31.75	ng/ul	0.00
81) Pyrene-d10	19.59	212	4613921	34.44	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.56	264	4899075	33.12	ng/ul	-0.01

Target Compounds Ovalue

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

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