

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP120520\
 Data File : BP004163.D
 Acq On : 05 Dec 2020 15:56
 Operator : CG/JU
 Sample : PB133233BL
 Misc : MDL-WATER-BL-02
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK233

Quant Time: Dec 07 02:55:07 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	388835	20.00	ng/ul	0.00
20) Naphthalene-d8	10.65	136	1613653	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.49	164	1021820	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.25	188	2266932	20.00	ng/ul	0.00
79) Chrysene-d12	21.34	240	2385272	20.00	ng/ul	0.00
88) Perylene-d12	23.71	264	2554412	20.00	ng/ul	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.35	96	64201	6.34	ng/uL	0.00
4) Pyridine-d5	3.74	84	887639	30.87	ng/ul	0.00
7) Phenol-d5	7.00	99	1041168	29.95	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.18	67	643727	32.03	ng/ul	0.00
11) 2-Chlorophenol-d4	7.38	132	804563	29.74	ng/ul	0.00
15) 4-Methylphenol-d8	8.55	113	792963	28.95	ng/ul	0.00
21) Nitrobenzene-d5	9.00	128	408585	31.06	ng/ul	0.00
24) 2-Nitrophenol-d4	9.72	143	358792	29.87	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.26	165	746465	28.24	ng/ul	0.00
31) 4-Chloroaniline-d4	10.77	131	1153595	29.74	ng/ul	0.00
46) Dimethylphthalate-d6	13.90	166	2677058	32.99	ng/ul	0.00
49) Acenaphthylene-d8	14.19	160	3015927	30.24	ng/ul	0.00
54) 4-Nitrophenol-d4	14.67	143	367649	27.68	ng/ul	0.00
60) Fluorene-d10	15.49	176	2231564	31.44	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.60	200	224009	24.32	ng/ul	0.00
73) Anthracene-d10	17.35	188	3583792	32.17	ng/ul	0.00
81) Pyrene-d10	19.59	212	4152214	33.09	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.56	264	4605306	33.56	ng/ul	0.00

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

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