

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120420\
 Data File : BM028032.D
 Acq On : 04 Dec 2020 16:05
 Operator : JU/CG
 Sample : PB133233BL
 Misc : MDL-WATER-BL-03
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK233

Quant Time: Dec 04 17:11:59 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120320.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 04 11:20:16 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	226711	20.00	ng/ul	0.00
20) Naphthalene-d8	11.13	136	979179	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.91	164	597133	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.63	188	1207904	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	1077555	20.00	ng/ul	0.00
88) Perylene-d12	24.14	264	1026183	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	34652	5.71	ng/uL	0.00
4) Pyridine-d5	3.93	84	489409	29.88	ng/ul	0.00
7) Phenol-d5	7.42	99	596294	31.60	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.60	67	385661	32.88	ng/ul	0.00
11) 2-Chlorophenol-d4	7.81	132	505400	32.15	ng/ul	0.00
15) 4-Methylphenol-d8	8.99	113	488482	32.82	ng/ul	0.00
21) Nitrobenzene-d5	9.47	128	243085	31.98	ng/ul	0.00
24) 2-Nitrophenol-d4	10.20	143	255618	33.65	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.73	165	466800	30.30	ng/ul	0.00
31) 4-Chloroaniline-d4	11.25	131	699952	33.21	ng/ul	0.00
46) Dimethylphthalate-d6	14.31	166	1535944	34.77	ng/ul	0.00
49) Acenaphthylene-d8	14.61	160	1730151	30.51	ng/ul	0.00
54) 4-Nitrophenol-d4	15.07	143	258668	30.36	ng/ul	0.00
60) Fluorene-d10	15.89	176	1267579	32.01	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	16.00	200	197944	28.23	ng/ul	0.00
73) Anthracene-d10	17.73	188	1927166	32.55	ng/ul	0.00
81) Pyrene-d10	19.98	212	2026028	35.82	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.99	264	1885559	34.51	ng/ul	0.00

Target Compounds Ovalue

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

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