

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120420\
 Data File : BM028030.D
 Acq On : 04 Dec 2020 14:52
 Operator : JU/CG
 Sample : PB133233BL
 Misc : MDL-WATER-BL-02
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK233

Quant Time: Dec 04 15:57:27 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	209634	20.00	ng/ul	0.00
20) Naphthalene-d8	11.13	136	833956	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.91	164	454390	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.63	188	872652	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	759906	20.00	ng/ul	0.00
88) Perylene-d12	24.14	264	757413	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.47	96	34062	6.07	ng/uL	0.00
4) Pyridine-d5	3.93	84	452942	29.91	ng/ul	0.00
7) Phenol-d5	7.42	99	539612	30.92	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.60	67	341108	31.45	ng/ul	0.00
11) 2-Chlorophenol-d4	7.81	132	460324	31.67	ng/ul	0.00
15) 4-Methylphenol-d8	8.99	113	425051	30.88	ng/ul	0.00
21) Nitrobenzene-d5	9.47	128	208855	32.26	ng/ul	0.00
24) 2-Nitrophenol-d4	10.20	143	215285	33.28	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.73	165	397537	30.30	ng/ul	0.00
31) 4-Chloroaniline-d4	11.25	131	575414	32.06	ng/ul	0.00
46) Dimethylphthalate-d6	14.30	166	1115201	33.17	ng/ul	0.00
49) Acenaphthylene-d8	14.61	160	1320307	30.60	ng/ul	0.00
54) 4-Nitrophenol-d4	15.07	143	183712	28.33	ng/ul	0.00
60) Fluorene-d10	15.89	176	936322	31.08	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.99	200	130419	25.75	ng/ul	0.00
73) Anthracene-d10	17.73	188	1392646	32.55	ng/ul	0.00
81) Pyrene-d10	19.98	212	1432999	35.92	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.99	264	1392112	34.52	ng/ul	0.00

Target Compounds Ovalue

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

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