

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120820\
 Data File : BM028077.D
 Acq On : 08 Dec 2020 18:09
 Operator : JU/CG
 Sample : PB133372BL
 Misc : MDL-SOIL-BL-01
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK372

Quant Time: Dec 09 03:45:20 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM120820.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 09 03:34:15 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.28	152	191062	20.00	ng/ul	0.00
20) Naphthalene-d8	11.12	136	773527	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.90	164	450142	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.62	188	922693	20.00	ng/ul	0.00
79) Chrysene-d12	21.73	240	889870	20.00	ng/ul	0.00
88) Perylene-d12	24.13	264	897533	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.46	96	29556	5.91	ng/uL	0.00
4) Pyridine-d5	3.92	84	396800	29.04	ng/ul	0.00
7) Phenol-d5	7.41	99	485543	29.77	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.59	67	310299	29.94	ng/ul	0.00
11) 2-Chlorophenol-d4	7.80	132	413988	30.49	ng/ul	0.00
15) 4-Methylphenol-d8	8.98	113	386065	29.00	ng/ul	0.00
21) Nitrobenzene-d5	9.46	128	193494	31.00	ng/ul	0.00
24) 2-Nitrophenol-d4	10.19	143	202419	31.49	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.72	165	368911	29.21	ng/ul	0.00
31) 4-Chloroaniline-d4	11.24	131	534310	30.33	ng/ul	0.00
46) Dimethylphthalate-d6	14.30	166	1168526	32.65	ng/ul	0.00
49) Acenaphthylene-d8	14.60	160	1305640	30.02	ng/ul	0.00
54) 4-Nitrophenol-d4	15.07	143	191021	28.33	ng/ul	0.00
60) Fluorene-d10	15.89	176	952165	30.58	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.99	200	149219	26.47	ng/ul	0.00
73) Anthracene-d10	17.72	188	1440170	31.36	ng/ul	0.00
81) Pyrene-d10	19.97	212	1597350	33.06	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.98	264	1637149	33.31	ng/ul	0.00

Target Compounds Ovalue

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

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