

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM031921\
 Data File : BM029173.D
 Acq On : 19 Mar 2021 16:39
 Operator : CG/JU
 Sample : PB134985BL
 Misc : SFAM-MDL-MES-BL01
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK985

Quant Time: Mar 19 17:14:35 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-BM031921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 19 12:41:26 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	75402	20.00	ng/ul	0.00
20) Naphthalene-d8	10.55	136	304072	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.39	164	200461	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.13	188	432014	20.00	ng/ul	0.00
79) Chrysene-d12	21.29	240	452255	20.00	ng/ul	0.00
88) Perylene-d12	23.53	264	486704	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.28	96	12030	6.27	ng/uL	0.00
4) Pyridine-d5	3.69	84	151786	30.27	ng/ul	0.00
7) Phenol-d5	6.93	99	196078	32.93	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.11	67	119394	34.69	ng/ul	0.00
11) 2-Chlorophenol-d4	7.30	132	165742	35.76	ng/ul	0.00
15) 4-Methylphenol-d8	8.46	113	156693	33.90	ng/ul	0.00
21) Nitrobenzene-d5	8.92	128	68055	35.19	ng/ul	0.00
24) 2-Nitrophenol-d4	9.64	143	70175	38.06	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.17	165	157381	32.58	ng/ul	0.00
31) 4-Chloroaniline-d4	10.69	131	223547	32.70	ng/ul	0.00
46) Dimethylphthalate-d6	13.81	166	539468	35.68	ng/ul	0.00
49) Acenaphthylene-d8	14.09	160	635767	35.57	ng/ul	0.00
54) 4-Nitrophenol-d4	14.57	143	71965	29.23	ng/ul	0.00
60) Fluorene-d10	15.39	176	448881	34.50	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.49	200	53108	26.08	ng/ul	0.00
73) Anthracene-d10	17.23	188	718956	34.85	ng/ul	0.00
81) Pyrene-d10	19.52	212	829649	39.31	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.39	264	927952	35.82	ng/ul	0.00

Target Compounds Ovalue

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

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