

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN031721\
 Data File : BN013969.D
 Acq On : 17 Mar 2021 14:22
 Operator : CG/JU
 Sample : PB134985BL
 Misc : SFAM-MDL-MES-BL-01
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK985

Quant Time: Mar 17 14:52:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN031721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 17 10:09:09 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	138309	20.00	ng/ul	0.00
20) Naphthalene-d8	10.49	136	555771	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.35	164	330767	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.09	188	673813	20.00	ng/ul	0.00
79) Chrysene-d12	21.27	240	681947	20.00	ng/ul	0.00
88) Perylene-d12	23.50	264	703820	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.21	96	24728	7.04	ng/uL	0.00
4) Pyridine-d5	3.61	84	314894	32.71	ng/ul	0.00
7) Phenol-d5	6.87	99	397331	33.91	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.05	67	254499	34.68	ng/ul	0.00
11) 2-Chlorophenol-d4	7.24	132	331048	37.44	ng/ul	0.00
15) 4-Methylphenol-d8	8.40	113	297294	33.06	ng/ul	0.00
21) Nitrobenzene-d5	8.86	128	142558	36.79	ng/ul	0.00
24) 2-Nitrophenol-d4	9.58	143	141272	38.51	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.10	165	280677	35.48	ng/ul	0.00
31) 4-Chloroaniline-d4	10.62	131	421325	34.75	ng/ul	0.00
46) Dimethylphthalate-d6	13.76	166	853131	37.19	ng/ul	0.00
49) Acenaphthylene-d8	14.04	160	1099107	39.28	ng/ul	0.00
54) 4-Nitrophenol-d4	14.54	143	126083	28.34	ng/ul	0.00
60) Fluorene-d10	15.34	176	753362	36.90	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.46	200	88590	26.30	ng/ul	0.00
73) Anthracene-d10	17.19	188	1181680	38.07	ng/ul	0.00
81) Pyrene-d10	19.49	212	1331621	38.66	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.36	264	1415096	39.92	ng/ul	0.00

Target Compounds Ovalue

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

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