

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP031821\
 Data File : BP004997.D
 Acq On : 18 Mar 2021 12:26
 Operator : CG/JU
 Sample : PB134984BL
 Misc : SFAM-MDL-MES-BL-01
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK985

Quant Time: Mar 18 13:11:45 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP031721.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 18 10:42:23 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	48988	20.00	ng/ul	0.00
20) Naphthalene-d8	10.52	136	192269	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.38	164	130356	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.13	188	296968	20.00	ng/ul	0.00
79) Chrysene-d12	21.23	240	326473	20.00	ng/ul	0.00
88) Perylene-d12	23.52	264	319196	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.23	96	6913	5.58	ng/uL	0.00
4) Pyridine-d5	3.64	84	89142	29.48	ng/ul	0.00
7) Phenol-d5	6.90	99	121397	30.52	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.06	67	61759	29.44	ng/ul	0.00
11) 2-Chlorophenol-d4	7.26	132	105331	32.93	ng/ul	0.00
15) 4-Methylphenol-d8	8.43	113	97691	30.96	ng/ul	0.00
21) Nitrobenzene-d5	8.88	128	48729	33.15	ng/ul	0.00
24) 2-Nitrophenol-d4	9.60	143	57201	34.42	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.14	165	106262	33.24	ng/ul	0.00
31) 4-Chloroaniline-d4	10.66	131	138741	31.92	ng/ul	0.00
46) Dimethylphthalate-d6	13.79	166	349464	35.10	ng/ul	0.00
49) Acenaphthylene-d8	14.07	160	413496	35.57	ng/ul	0.00
54) 4-Nitrophenol-d4	14.59	143	47369	27.05	ng/ul	0.00
60) Fluorene-d10	15.37	176	294473	34.18	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.50	200	59924	28.90	ng/ul	0.00
73) Anthracene-d10	17.23	188	480164	34.37	ng/ul	0.00
81) Pyrene-d10	19.48	212	567289	34.97	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.37	264	628164	36.59	ng/ul	0.00

Target Compounds Ovalue

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

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