

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011421\
 Data File : BG047914.D
 Acq On : 14 Jan 2021 10:11
 Operator : CG/JU
 Sample : PB133977BL
 Misc : PBBL-SFAM-MDL-W-01
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK977

Quant Time: Jan 14 11:30:22 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG011321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 13 15:41:08 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.15	152	78660	20.00	ng/ul	0.00
20) Naphthalene-d8	10.97	136	303710	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.78	164	207479	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.52	188	512899	20.00	ng/ul	0.00
79) Chrysene-d12	21.80	240	588580	20.00	ng/ul	0.00
88) Perylene-d12	25.11	264	534280	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.57	96	15757	7.02	ng/uL	0.00
4) Pyridine-d5	3.98	84	234938	39.39	ng/ul	0.00
7) Phenol-d5	7.29	99	266992	39.38	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.47	67	169603	39.36	ng/ul	0.00
11) 2-Chlorophenol-d4	7.68	132	204392	39.48	ng/ul	0.00
15) 4-Methylphenol-d8	8.84	113	207735	38.78	ng/ul	0.00
21) Nitrobenzene-d5	9.31	128	88833	46.81	ng/ul	0.00
24) 2-Nitrophenol-d4	10.04	143	91081	45.35	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.58	165	210996	39.05	ng/ul	0.00
31) 4-Chloroaniline-d4	11.09	131	286487	39.73	ng/ul	0.00
46) Dimethylphthalate-d6	14.18	166	718408	41.99	ng/ul	0.00
49) Acenaphthylene-d8	14.47	160	774632	38.30	ng/ul	0.00
54) 4-Nitrophenol-d4	14.94	143	97711	38.79	ng/ul	0.00
60) Fluorene-d10	15.76	176	605308	39.64	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.87	200	83160	36.19	ng/ul	0.00
73) Anthracene-d10	17.62	188	978798	39.83	ng/ul	0.00
81) Pyrene-d10	19.89	212	1186864	37.06	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.89	264	1253674	42.35	ng/ul	0.01

Target Compounds Ovalue

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

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