

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011421\
 Data File : BG047916.D
 Acq On : 14 Jan 2021 13:13
 Operator : CG/JU
 Sample : PB134057BL
 Misc : PBBL-SFAM-MDL-S-01
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK057

Quant Time: Jan 14 13:57:20 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	8.16	152	77382	20.00	ng/ul	0.00
20) Naphthalene-d8	10.97	136	298880	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.78	164	205250	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.52	188	490668	20.00	ng/ul	0.00
79) Chrysene-d12	21.80	240	546933	20.00	ng/ul	0.00
88) Perylene-d12	25.12	264	498503	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.56	96	15592	7.06	ng/uL	0.00
4) Pyridine-d5	3.98	84	219954	37.49	ng/ul	0.00
7) Phenol-d5	7.29	99	258852	38.81	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.47	67	170034	40.11	ng/ul	0.00
11) 2-Chlorophenol-d4	7.68	132	197746	38.83	ng/ul	0.00
15) 4-Methylphenol-d8	8.85	113	199571	37.87	ng/ul	0.00
21) Nitrobenzene-d5	9.32	128	88410	47.34	ng/ul	0.00
24) 2-Nitrophenol-d4	10.04	143	90794	45.94	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.58	165	205060	38.56	ng/ul	0.00
31) 4-Chloroaniline-d4	11.10	131	277068	39.04	ng/ul	0.00
46) Dimethylphthalate-d6	14.17	166	683505	40.38	ng/ul	0.00
49) Acenaphthylene-d8	14.47	160	752871	37.63	ng/ul	0.00
54) 4-Nitrophenol-d4	14.94	143	94511	37.92	ng/ul	0.00
60) Fluorene-d10	15.77	176	579367	38.36	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.87	200	78149	35.55	ng/ul	0.00
73) Anthracene-d10	17.62	188	934445	39.74	ng/ul	0.00
81) Pyrene-d10	19.90	212	1123440	37.75	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.89	264	1174958	42.54	ng/ul	0.00

Target Compounds Ovalue

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

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