

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG020421\
 Data File : BG048112.D
 Acq On : 4 Feb 2021 12:42
 Operator : CG/JU
 Sample : PB134021BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK021

Quant Time: Feb 04 14:00:41 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG012921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Feb 04 12:55:42 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	50198	20.00	ng/ul	0.00
20) Naphthalene-d8	10.92	136	195719	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.74	164	143121	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.47	188	359520	20.00	ng/ul	0.00
79) Chrysene-d12	21.75	240	399525	20.00	ng/ul	0.00
88) Perylene-d12	25.02	264	378401	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.52	96	9075	6.80	ng/uL	0.00
4) Pyridine-d5	3.93	84	141619	36.33	ng/ul	0.00
7) Phenol-d5	7.26	99	178414	38.12	ng/ul	0.00
9) Bis-(2-Chloroethyl)ether-d	7.43	67	110107	39.05	ng/ul	0.00
11) 2-Chlorophenol-d4	7.64	132	128678	37.78	ng/ul	0.00
15) 4-Methylphenol-d8	8.80	113	138388	37.46	ng/ul	0.00
21) Nitrobenzene-d5	9.27	128	66983	40.70	ng/ul	0.00
24) 2-Nitrophenol-d4	9.99	143	79611	39.36	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.53	165	139572	36.47	ng/ul	0.00
31) 4-Chloroaniline-d4	11.05	131	177688	36.25	ng/ul	0.00
46) Dimethylphthalate-d6	14.14	166	479316	40.10	ng/ul	0.00
49) Acenaphthylene-d8	14.43	160	506116	36.79	ng/ul	0.00
54) 4-Nitrophenol-d4	14.91	143	79749	37.20	ng/ul	0.00
60) Fluorene-d10	15.72	176	407998	38.09	ng/ul	0.00
65) 4,6-Dinitro-2-methylphenol	15.83	200	92133	35.83	ng/ul	0.00
73) Anthracene-d10	17.57	188	665703	38.35	ng/ul	0.00
81) Pyrene-d10	19.85	212	846751	38.02	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.79	264	878996	42.49	ng/ul	0.00

Target Compounds Ovalue

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

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