

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070524\
 Data File : BG061913.D
 Acq On : 6 Jul 2024 2:36
 Operator : MA/JU
 Sample : P2982-18
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A4BB9

Quant Time: Jul 06 03:51:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jul 03 02:38:46 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.094	152	33444	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.908	136	170389	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.715	164	127893	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.459	188	298759	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.719	240	273187	20.000	ng/ul	0.00
88) Perylene-d12	24.962	264	324196	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.487	96	3114	3.499	ng/uL	0.00
4) Pyridine-d5	3.904	84	44516	16.268	ng/ul	0.00
7) Phenol-d5	7.247	99	87218	23.830	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.418	67	52343	22.584	ng/ul	0.00
11) 2-Chlorophenol-d4	7.624	132	63515	25.299	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	72650	25.211	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	32531	25.104	ng/ul	0.00
24) 2-Nitrophenol-d4	9.985	143	40413	27.309	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.520	165	77513	27.156	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.043	131	100984	24.368	ng/ul	0.00
46) Dimethylphthalate-d6	14.116	166	274728	26.128	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	306401	27.869	ng/ul	0.00
54) 4-Nitrophenol-d4	14.903	143	40125	23.896	ng/ul	-0.01
60) Fluorene-d10	15.702	176	236773	26.749	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.814	200	36311	23.142	ng/ul	-0.01
73) Anthracene-d10	17.553	188	375282	26.905	ng/ul	0.00
81) Pyrene-d10	19.833	212	432483	26.915	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.733	264	452299	28.131	ng/ul	0.00

Target Compounds Qvalue

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

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