

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100923\
 Data File : BG059417.D
 Acq On : 10 Oct 2023 6:09
 Operator : MA/JU
 Sample : PB156059BL
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK059

Quant Time: Oct 10 06:48:55 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG100723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 07 22:45:22 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.223	152	39875	20.000	ng/u1	0.00
20) Naphthalene-d8	11.067	136	186455	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.856	164	129691	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.606	188	304923	20.000	ng/u1	0.00
79) Chrysene-d12	21.907	240	268886	20.000	ng/u1	0.00
88) Perylene-d12	25.391	264	323388	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.523	96	9520	8.512	ng/uL	0.00
4) Pyridine-d5	3.963	84	123657	38.414	ng/u1	0.00
7) Phenol-d5	7.353	99	168372	38.241	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	93319	37.499	ng/u1	0.00
11) 2-Chlorophenol-d4	7.741	132	120133	39.013	ng/u1	0.00
15) 4-Methylphenol-d8	8.916	113	126075	36.752	ng/u1	0.00
21) Nitrobenzene-d5	9.427	128	63736	40.155	ng/u1	0.00
24) 2-Nitrophenol-d4	10.144	143	74862	43.577	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.673	165	131414	42.381	ng/u1	0.00
31) 4-Chloroaniline-d4	11.213	131	192830	39.150	ng/u1	0.00
46) Dimethylphthalate-d6	14.257	166	446401	42.014	ng/u1	0.00
49) Acenaphthylene-d8	14.563	160	513731	40.630	ng/u1	0.00
54) 4-Nitrophenol-d4	15.050	143	70096	35.205	ng/u1	0.00
60) Fluorene-d10	15.849	176	406600	40.091	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.961	200	62391	30.991	ng/u1	0.00
73) Anthracene-d10	17.706	188	658842	44.046	ng/u1	0.00
81) Pyrene-d10	19.986	212	728316	44.712	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.150	264	759162	44.611	ng/u1	0.00

Target Compounds Qvalue

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

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