

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100723\
 Data File : BG059362.D
 Acq On : 8 Oct 2023 5:06
 Operator : MA/JU
 Sample : PB155977BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK977

Quant Time: Oct 08 05:43:41 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	8.230	152	43148	20.000	ng/u1	0.00
20) Naphthalene-d8	11.067	136	205720	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.863	164	141732	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.613	188	337154	20.000	ng/u1	0.00
79) Chrysene-d12	21.914	240	303302	20.000	ng/u1	0.00
88) Perylene-d12	25.404	264	352056	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.523	96	9636	7.963	ng/uL	0.00
4) Pyridine-d5	3.958	84	146775	42.137	ng/u1	0.00
7) Phenol-d5	7.354	99	190725	40.032	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.548	67	111196	41.293	ng/u1	0.00
11) 2-Chlorophenol-d4	7.748	132	142964	42.905	ng/u1	0.00
15) 4-Methylphenol-d8	8.917	113	143430	38.639	ng/u1	0.00
21) Nitrobenzene-d5	9.428	128	76356	43.601	ng/u1	0.00
24) 2-Nitrophenol-d4	10.145	143	79736	42.067	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.674	165	146741	42.892	ng/u1	0.00
31) 4-Chloroaniline-d4	11.214	131	221441	40.749	ng/u1	0.00
46) Dimethylphthalate-d6	14.258	166	518789	44.679	ng/u1	0.00
49) Acenaphthylene-d8	14.563	160	603563	43.679	ng/u1	0.00
54) 4-Nitrophenol-d4	15.051	143	87396	40.164	ng/u1	0.00
60) Fluorene-d10	15.850	176	467418	42.172	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.962	200	86759	38.975	ng/u1	0.00
73) Anthracene-d10	17.713	188	754788	45.637	ng/u1	0.00
81) Pyrene-d10	19.986	212	830332	45.191	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.157	264	847731	45.759	ng/u1	0.00

Target Compounds	Qvalue

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

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