

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG100723\
 Data File : BG059376.D
 Acq On : 8 Oct 2023 15:12
 Operator : MA/JU
 Sample : PB156048BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK048

Quant Time: Oct 09 01:40:28 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	8.229	152	50853	20.000	ng/u1	0.00
20) Naphthalene-d8	11.067	136	239521	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.862	164	162309	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.612	188	379379	20.000	ng/u1	0.00
79) Chrysene-d12	21.913	240	345561	20.000	ng/u1	0.00
88) Perylene-d12	25.403	264	409230	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.523	96	11478	8.048	ng/uL	0.00
4) Pyridine-d5	3.963	84	156140	38.034	ng/u1	0.00
7) Phenol-d5	7.354	99	204985	36.506	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.547	67	118530	37.347	ng/u1	0.00
11) 2-Chlorophenol-d4	7.747	132	147580	37.580	ng/u1	0.00
15) 4-Methylphenol-d8	8.922	113	150567	34.416	ng/u1	0.00
21) Nitrobenzene-d5	9.422	128	78748	38.622	ng/u1	0.00
24) 2-Nitrophenol-d4	10.144	143	88672	40.180	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.673	165	154041	38.672	ng/u1	0.00
31) 4-Chloroaniline-d4	11.214	131	237906	37.600	ng/u1	0.00
46) Dimethylphthalate-d6	14.257	166	543191	40.850	ng/u1	0.00
49) Acenaphthylene-d8	14.563	160	624992	39.496	ng/u1	0.00
54) 4-Nitrophenol-d4	15.050	143	91476	36.710	ng/u1	0.00
60) Fluorene-d10	15.849	176	496078	39.084	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.961	200	89208	35.615	ng/u1	0.00
73) Anthracene-d10	17.712	188	768893	41.315	ng/u1	0.00
81) Pyrene-d10	19.986	212	860978	41.128	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.156	264	893247	41.480	ng/u1	0.00

Target Compounds	Qvalue

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

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