

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG071123\  
 Data File : BG058154.D  
 Acq On : 11 Jul 2023 10:50  
 Operator : CG/JU  
 Sample : PB153937BL  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :

Quant Time: Jul 11 22:53:58 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG070123.MA.M  
 Quant Title : SVOA CALIBRATION  
 QLast Update : Mon Jul 10 16:52:40 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.971	152	47967	20.000	ng/ul	0.00
20) Naphthalene-d8	10.780	136	215021	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.605	164	147191	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.343	188	345618	20.000	ng/ul	0.00
79) Chrysene-d12	21.602	240	305292	20.000	ng/ul	0.00
88) Perylene-d12	24.799	264	357955	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.400	96	7049	5.146	ng/uL	0.00
4) Pyridine-d5	3.811	84	110197	26.498	ng/ul	0.00
7) Phenol-d5	7.131	99	143019	29.196	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.296	67	85792	29.062	ng/ul	0.00
11) 2-Chlorophenol-d4	7.501	132	98702	29.214	ng/ul	0.00
15) 4-Methylphenol-d8	8.676	113	108690	29.405	ng/ul	0.00
21) Nitrobenzene-d5	9.135	128	52551	29.986	ng/ul	0.00
24) 2-Nitrophenol-d4	9.857	143	59200	32.141	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.398	165	105009	29.719	ng/ul	0.00
31) 4-Chloroaniline-d4	10.909	131	150468	28.088	ng/ul	0.00
46) Dimethylphthalate-d6	14.005	166	327601	28.208	ng/ul	0.00
49) Acenaphthylene-d8	14.299	160	384117	30.310	ng/ul	0.00
54) 4-Nitrophenol-d4	14.799	143	52036	25.361	ng/ul	0.00
60) Fluorene-d10	15.592	176	299539	30.370	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.709	200	40202	21.421	ng/ul	0.00
73) Anthracene-d10	17.443	188	488650	30.469	ng/ul	0.00
81) Pyrene-d10	19.728	212	556545	28.945	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.581	264	596452	32.706	ng/ul	0.01

Target Compounds Qvalue

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

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