

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG072223\
 Data File : BG058401.D
 Acq On : 22 Jul 2023 14:00
 Operator : CG/JU
 Sample : PB154168BL
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Jul 23 01:31:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 18 02:19:39 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.896	152	49834	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	217968	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.535	164	147843	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.279	188	332808	20.000	ng/ul	0.00
79) Chrysene-d12	21.527	240	284462	20.000	ng/ul	0.00
88) Perylene-d12	24.670	264	348467	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.337	96	8890	6.566	ng/uL	0.00
4) Pyridine-d5	3.748	84	139948	34.824	ng/ul	0.00
7) Phenol-d5	7.056	99	162476	34.369	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.220	67	96145	32.539	ng/ul	0.00
11) 2-Chlorophenol-d4	7.426	132	116571	35.227	ng/ul	0.00
15) 4-Methylphenol-d8	8.601	113	120088	33.303	ng/ul	0.00
21) Nitrobenzene-d5	9.053	128	58801	35.990	ng/ul	0.00
24) 2-Nitrophenol-d4	9.776	143	64868	36.368	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.317	165	116785	34.156	ng/ul	0.00
31) 4-Chloroaniline-d4	10.834	131	159774	32.234	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	367116	32.202	ng/ul	0.00
49) Acenaphthylene-d8	14.230	160	422987	35.404	ng/ul	0.00
54) 4-Nitrophenol-d4	14.741	143	43566	25.304	ng/ul	0.00
60) Fluorene-d10	15.528	176	325875	34.525	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.646	200	27340	16.124	ng/ul	0.00
73) Anthracene-d10	17.379	188	520434	36.191	ng/ul	0.00
81) Pyrene-d10	19.670	212	600106	35.718	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.453	264	631266	37.481	ng/ul	0.00

Target Compounds Qvalue

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

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