

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071123\
 Data File : BG058165.D
 Acq On : 11 Jul 2023 19:04
 Operator : CG/JU
 Sample : PB153960BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :

Quant Time: Jul 11 22:57:52 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	7.975	152	48088	20.000	ng/ul	0.00
20) Naphthalene-d8	10.778	136	225239	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.603	164	157916	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.341	188	368953	20.000	ng/ul	0.00
79) Chrysene-d12	21.601	240	328117	20.000	ng/ul	0.00
88) Perylene-d12	24.791	264	384703	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.404	96	6549	4.769	ng/uL	0.00
4) Pyridine-d5	3.816	84	111404	26.721	ng/ul	0.00
7) Phenol-d5	7.135	99	145695	29.667	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.294	67	85660	28.944	ng/ul	0.00
11) 2-Chlorophenol-d4	7.505	132	101568	29.987	ng/ul	0.00
15) 4-Methylphenol-d8	8.675	113	112285	30.301	ng/ul	0.00
21) Nitrobenzene-d5	9.133	128	54940	29.927	ng/ul	0.00
24) 2-Nitrophenol-d4	9.856	143	62608	32.449	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.396	165	109813	29.669	ng/ul	0.00
31) 4-Chloroaniline-d4	10.907	131	158498	28.244	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	356327	28.598	ng/ul	0.00
49) Acenaphthylene-d8	14.297	160	409112	30.089	ng/ul	0.00
54) 4-Nitrophenol-d4	14.797	143	52421	23.813	ng/ul	0.00
60) Fluorene-d10	15.590	176	318242	30.075	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.708	200	45110	22.516	ng/ul	0.00
73) Anthracene-d10	17.441	188	510055	29.792	ng/ul	0.00
81) Pyrene-d10	19.726	212	593830	28.736	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.574	264	626348	31.957	ng/ul	0.00

Target Compounds Qvalue

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

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