

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058991.D
 Acq On : 11 Sep 2023 20:57
 Operator : MA/JU
 Sample : PB155374BL
 Misc :
 ALS Vial : 91 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK374

Quant Time: Sep 12 01:50:30 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 00:07:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.325	152	54649	20.000	ng/u1	0.00
20) Naphthalene-d8	11.169	136	277181	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.947	164	223651	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.697	188	584879	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.015	240	551202	20.000	ng/u1	-0.01
88) Perylene-d12	25.587	264	619230	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.637	96	11949	9.294	ng/uL	0.00
4) Pyridine-d5	4.066	84	172711	42.429	ng/u1	0.00
7) Phenol-d5	7.432	99	222420	40.360	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.644	67	129173	39.974	ng/u1	0.00
11) 2-Chlorophenol-d4	7.838	132	147048	41.604	ng/u1	0.00
15) 4-Methylphenol-d8	9.001	113	180001	42.278	ng/u1	0.00
21) Nitrobenzene-d5	9.518	128	77913	38.752	ng/u1	0.00
24) 2-Nitrophenol-d4	10.235	143	92879	42.100	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.758	165	181136	38.633	ng/u1	0.00
31) 4-Chloroaniline-d4	11.310	131	238586	37.840	ng/u1	0.00
46) Dimethylphthalate-d6	14.342	166	649525	37.440	ng/u1	0.00
49) Acenaphthylene-d8	14.647	160	681585	35.281	ng/u1	0.00
54) 4-Nitrophenol-d4	15.112	143	115245	39.992	ng/u1	0.00
60) Fluorene-d10	15.934	176	579110	37.615	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.034	200	99152	34.536	ng/u1	0.00
73) Anthracene-d10	17.797	188	956171	36.827	ng/u1	0.00
81) Pyrene-d10	20.065	212	1118904	38.321	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.335	264	1149307	37.862	ng/u1	0.00

Target Compounds Qvalue

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

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