

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090221\
 Data File : BG050131.D
 Acq On : 3 Sep 2021 14:51
 Operator : CG/JU
 Sample : PB138772BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK772

Quant Time: Sep 03 15:28:54 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG081921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Aug 30 14:59:55 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.950	152	25978	20.000	ng/ul	0.00
20) Naphthalene-d8	10.770	136	109250	20.000	ng/ul #	0.00
38) Acenaphthene-d10	14.612	164	72650	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.367	188	158377	20.000	ng/ul	0.00
79) Chrysene-d12	21.644	240	164273	20.000	ng/ul	0.00
88) Perylene-d12	24.863	264	160176	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	4673	9.276	ng/uL	0.00
4) Pyridine-d5	3.733	84	58140	38.315	ng/ul	0.00
7) Phenol-d5	7.122	99	86355	39.146	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	49204	40.149	ng/ul	0.00
11) 2-Chlorophenol-d4	7.480	132	66521	40.301	ng/ul	0.00
15) 4-Methylphenol-d8	8.673	113	69615	40.071	ng/ul	0.00
21) Nitrobenzene-d5	9.125	128	33853	39.969	ng/ul	0.00
24) 2-Nitrophenol-d4	9.848	143	41083	40.735	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.400	165	66203	36.803	ng/ul	0.00
31) 4-Chloroaniline-d4	10.911	131	90314	36.886	ng/ul	0.00
46) Dimethylphthalate-d6	14.013	166	220154	39.596	ng/ul	0.00
49) Acenaphthylene-d8	14.307	160	272220	41.757	ng/ul	0.00
54) 4-Nitrophenol-d4	14.841	143	27494	34.178	ng/ul	0.00
60) Fluorene-d10	15.611	176	183485	38.923	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.752	200	30433	29.868	ng/ul	0.00
73) Anthracene-d10	17.467	188	302704	42.563	ng/ul	0.00
81) Pyrene-d10	19.758	212	356459	40.155	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.640	264	349112	41.083	ng/ul	0.00

Target Compounds Qvalue

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

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