

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081822\
 Data File : BG054655.D
 Acq On : 18 Aug 2022 17:01
 Operator : CG/JU
 Sample : PB146659BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK659

Quant Time: Aug 18 18:59:27 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 18 03:22:26 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.165	152	48656	20.000	ng/u1	0.00
20) Naphthalene-d8	10.991	136	198506	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.792	164	131379	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.542	188	296977	20.000	ng/u1	0.00
79) Chrysene-d12	21.837	240	294188	20.000	ng/u1	0.00
88) Perylene-d12	25.215	264	293908	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.511	96	6429	5.227	ng/uL	0.00
4) Pyridine-d5	3.940	84	75916	20.711	ng/u1	0.00
7) Phenol-d5	7.295	99	104579	25.280	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.477	67	68683	25.947	ng/u1	0.00
11) 2-Chlorophenol-d4	7.689	132	77491	26.486	ng/u1	0.00
15) 4-Methylphenol-d8	8.858	113	81984	26.235	ng/u1	0.00
21) Nitrobenzene-d5	9.334	128	37103	29.777	ng/u1	0.00
24) 2-Nitrophenol-d4	10.063	143	33160	29.725	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.597	165	77541	26.514	ng/u1	0.00
31) 4-Chloroaniline-d4	11.114	131	116358	27.715	ng/u1	0.00
46) Dimethylphthalate-d6	14.193	166	269253	29.827	ng/u1	0.00
49) Acenaphthylene-d8	14.493	160	320745	29.676	ng/u1	0.00
54) 4-Nitrophenol-d4	14.957	143	38330	26.868	ng/u1	0.00
60) Fluorene-d10	15.785	176	232138	28.359	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.885	200	24356	24.083	ng/u1	0.00
73) Anthracene-d10	17.636	188	383959	29.725	ng/u1	0.00
81) Pyrene-d10	19.922	212	452411	29.646	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.980	264	436836	30.266	ng/u1	0.00

Target Compounds Qvalue

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

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