

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110821\
 Data File : BG050881.D
 Acq On : 8 Nov 2021 15:31
 Operator : CG/JU
 Sample : M4492-04
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG1S1

Quant Time: Nov 09 11:50:02 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 02 14:49:05 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.232	152	33355	20.000	ng/u1	0.00
20) Naphthalene-d8	11.058	136	148147	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.853	164	102697	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.597	188	225946	20.000	ng/u1	0.00
79) Chrysene-d12	21.892	240	194632	20.000	ng/u1	#-0.01
88) Perylene-d12	25.294	264	192117	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.561	96	18844	18.234	ng/uL	-0.03
4) Pyridine-d5	4.019	84	25600	8.280	ng/u1	0.00
7) Phenol-d5	7.374	99	27355	7.687	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	72866	31.700	ng/u1	0.00
11) 2-Chlorophenol-d4	7.756	132	62375	25.293	ng/u1	0.00
15) 4-Methylphenol-d8	8.925	113	45387	16.202	ng/u1	0.00
21) Nitrobenzene-d5	9.401	128	40448	32.128	ng/u1	0.00
24) 2-Nitrophenol-d4	10.130	143	41719	29.802	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.670	165	64223	27.235	ng/u1	0.00
31) 4-Chloroaniline-d4	11.187	131	91547	25.636	ng/u1	0.00
46) Dimethylphthalate-d6	14.248	166	270190	34.389	ng/u1	0.00
49) Acenaphthylene-d8	14.554	160	325808	33.284	ng/u1	0.00
54) 4-Nitrophenol-d4	15.041	143	9557	6.709	ng/u1	0.00
60) Fluorene-d10	15.840	176	242612	34.857	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.958	200	40141	29.303	ng/u1	0.00
73) Anthracene-d10	17.697	188	433020	40.536	ng/u1	0.00
81) Pyrene-d10	19.971	212	491004	39.058	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.059	264	408889	38.500	ng/u1	-0.01

Target Compounds Qvalue

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

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