

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG050972.D
 Acq On : 11 Nov 2021 17:42
 Operator : CG/JU
 Sample : M4522-10
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG3A0

Quant Time: Nov 12 01:45:27 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 11 12:40:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.230	152	30580	20.000	ng/u1	0.00
20) Naphthalene-d8	11.056	136	136062	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.851	164	94566	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.601	188	224173	20.000	ng/u1	0.00
79) Chrysene-d12	21.890	240	205756	20.000	ng/u1	0.00
88) Perylene-d12	25.298	264	200425	20.000	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	4228	4.462	ng/uL	0.00
4) Pyridine-d5	4.035	84	10397	3.668	ng/u1	0.02
7) Phenol-d5	7.372	99	19981	6.125	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	55567	26.368	ng/u1	0.00
11) 2-Chlorophenol-d4	7.760	132	49118	21.725	ng/u1	0.00
15) 4-Methylphenol-d8	8.929	113	35461	13.807	ng/u1	0.00
21) Nitrobenzene-d5	9.405	128	32842	28.404	ng/u1	0.00
24) 2-Nitrophenol-d4	10.127	143	35482	27.598	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.668	165	54678	25.247	ng/u1	0.00
31) 4-Chloroaniline-d4	11.191	131	68705	20.949	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	229390	31.706	ng/u1	0.00
49) Acenaphthylene-d8	14.552	160	264716	29.368	ng/u1	0.00
54) 4-Nitrophenol-d4	15.045	143	9344	7.123	ng/u1	0.01
60) Fluorene-d10	15.838	176	202473	31.591	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.956	200	41199	30.313	ng/u1	0.00
73) Anthracene-d10	17.695	188	376574	35.530	ng/u1	0.00
81) Pyrene-d10	19.975	212	458107	34.471	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.057	264	386313	34.866	ng/u1	0.00

Target Compounds Qvalue

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

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