

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110321\
 Data File : BG050872.D
 Acq On : 4 Nov 2021 7:53
 Operator : CG/JU
 Sample : M4412-17
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG371

Quant Time: Nov 07 08:33:48 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	8.230	152	30939	20.000	ng/u1	0.00
20) Naphthalene-d8	11.050	136	135056	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.851	164	98002	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.595	188	228677	20.000	ng/u1	-0.01
79) Chrysene-d12	21.890	240	203921	20.000	ng/u1	-0.02
88) Perylene-d12	25.286	264	198047	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.588	96	4614	4.813	ng/uL	0.00
4) Pyridine-d5	4.017	84	19844	6.920	ng/u1	0.00
7) Phenol-d5	7.372	99	24083	7.296	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	64567	30.283	ng/u1	0.00
11) 2-Chlorophenol-d4	7.754	132	56202	24.570	ng/u1	-0.01
15) 4-Methylphenol-d8	8.923	113	41919	16.132	ng/u1	-0.01
21) Nitrobenzene-d5	9.399	128	37391	32.579	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.127	143	38535	30.196	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.668	165	60778	28.273	ng/u1	0.00
31) 4-Chloroaniline-d4	11.185	131	75715	23.258	ng/u1	0.00
46) Dimethylphthalate-d6	14.246	166	263298	35.117	ng/u1	0.00
49) Acenaphthylene-d8	14.546	160	302000	32.329	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.045	143	9093	6.689	ng/u1	0.00
60) Fluorene-d10	15.838	176	229785	34.596	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.950	200	39799	28.706	ng/u1	-0.01
73) Anthracene-d10	17.695	188	400256	37.021	ng/u1	-0.01
81) Pyrene-d10	19.969	212	472316	35.860	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.057	264	386014	35.258	ng/u1	-0.02

Target Compounds Qvalue

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

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