

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG112321\
 Data File : BG051221.D
 Acq On : 24 Nov 2021 20:58
 Operator : CG/JU
 Sample : M4752-03
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG1W9

Quant Time: Nov 25 00:10:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 24 06:04:50 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.197	152	36990	20.000	ng/ul	0.00
20) Naphthalene-d8	11.029	136	171856	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.830	164	114712	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.580	188	229474	20.000	ng/ul	0.00
79) Chrysene-d12	21.881	240	191301	20.000	ng/ul	0.00
88) Perylene-d12	25.289	264	193663	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.544	96	4189	3.935	ng/uL	0.00
4) Pyridine-d5	3.972	84	18280	5.852	ng/ul	0.00
7) Phenol-d5	7.363	99	27329	7.475	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.515	67	66676	29.039	ng/ul	0.00
11) 2-Chlorophenol-d4	7.733	132	64076	24.340	ng/ul	0.00
15) 4-Methylphenol-d8	8.914	113	48742	16.522	ng/ul	0.00
21) Nitrobenzene-d5	9.378	128	43196	29.776	ng/ul	0.00
24) 2-Nitrophenol-d4	10.101	143	48493	29.633	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.653	165	78415	28.242	ng/ul	0.00
31) 4-Chloroaniline-d4	11.164	131	80712	19.867	ng/ul	0.00
46) Dimethylphthalate-d6	14.225	166	274665	31.119	ng/ul	0.00
49) Acenaphthylene-d8	14.531	160	339193	30.476	ng/ul	0.00
54) 4-Nitrophenol-d4	15.060	143	10378	7.264	ng/ul	0.01
60) Fluorene-d10	15.823	176	241484	30.382	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.953	200	36471	25.756	ng/ul	0.00
73) Anthracene-d10	17.680	188	362753	33.053	ng/ul	0.00
81) Pyrene-d10	19.960	212	416808	36.009	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.054	264	366611	35.445	ng/ul	0.01
Target Compounds						
8) Phenol	7.392	94	4004	1.057	ng/ul	96

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

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