

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
 Data File : BG049292.D
 Acq On : 11 Jul 2021 9:27
 Operator : CG/JU
 Sample : M2914-11
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBK29

Quant Time: Jul 12 01:04:14 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jul 09 03:15:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.074	152	28062	20.000	ng/ul	# 0.00
20) Naphthalene-d8	10.894	136	113145	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.719	164	81409	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	189923	20.000	ng/ul	0.00
79) Chrysene-d12	21.746	240	199805	20.000	ng/ul	0.00
88) Perylene-d12	25.042	264	209160	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.462	96	2994	5.019	ng/uL	0.00
4) Pyridine-d5	3.897	84	18935	10.794	ng/ul	0.00
7) Phenol-d5	7.222	99	19186	7.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.387	67	39415	27.840	ng/ul	0.00
11) 2-Chlorophenol-d4	7.598	132	46477	25.762	ng/ul	0.00
15) 4-Methylphenol-d8	8.773	113	38631	19.752	ng/ul	0.00
21) Nitrobenzene-d5	9.237	128	27987	32.234	ng/ul	0.00
24) 2-Nitrophenol-d4	9.966	143	31131	31.235	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.507	165	69776	35.878	ng/ul	0.00
31) 4-Chloroaniline-d4	11.024	131	70978	28.006	ng/ul	0.00
46) Dimethylphthalate-d6	14.114	166	261298	41.735	ng/ul	0.00
49) Acenaphthylene-d8	14.414	160	288406	39.261	ng/ul	0.00
54) 4-Nitrophenol-d4	14.913	143	7613	7.346	ng/ul	0.00
60) Fluorene-d10	15.706	176	209584	39.538	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.818	200	22899	18.969	ng/ul	0.00
73) Anthracene-d10	17.569	188	368913	42.664	ng/ul	0.00
81) Pyrene-d10	19.854	212	434594	42.437	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.819	264	465089	42.775	ng/ul	0.00

Target Compounds Qvalue

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

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