

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG070921\
 Data File : BG049287.D
 Acq On : 11 Jul 2021 6:03
 Operator : CG/JU
 Sample : M2914-13
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBK42

Quant Time: Jul 11 06:41:07 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG070921.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Jul 11 04:28:44 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.075	152	25750	20.000	ng/ul	0.00
20) Naphthalene-d8	10.889	136	112532	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.714	164	81024	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.464	188	190663	20.000	ng/ul	0.00
79) Chrysene-d12	21.747	240	194874	20.000	ng/ul	0.00
88) Perylene-d12	25.049	264	206648	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.468	96	3581	6.542	ng/uL	0.00
4) Pyridine-d5	3.891	84	19446	12.081	ng/ul	0.00
7) Phenol-d5	7.223	99	21533	9.614	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.393	67	44935	34.589	ng/ul	0.00
11) 2-Chlorophenol-d4	7.599	132	49981	30.192	ng/ul	0.00
15) 4-Methylphenol-d8	8.774	113	38094	21.227	ng/ul	0.00
21) Nitrobenzene-d5	9.238	128	30132	34.894	ng/ul	0.00
24) 2-Nitrophenol-d4	9.966	143	35701	36.015	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.507	165	69300	35.828	ng/ul	0.00
31) 4-Chloroaniline-d4	11.024	131	74433	29.529	ng/ul	0.00
46) Dimethylphthalate-d6	14.115	166	249128	39.981	ng/ul	0.00
49) Acenaphthylene-d8	14.408	160	283495	38.775	ng/ul	0.00
54) 4-Nitrophenol-d4	14.908	143	8967	8.694	ng/ul	0.00
60) Fluorene-d10	15.707	176	208740	39.566	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.818	200	32401	26.736	ng/ul	0.00
73) Anthracene-d10	17.563	188	353998	40.780	ng/ul	0.00
81) Pyrene-d10	19.849	212	405885	40.637	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.820	264	439084	40.875	ng/ul	0.00

Target Compounds Qvalue

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

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