

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062892.D
 Acq On : 17 Sep 2024 00:05
 Operator : JU/RC
 Sample : SP6631
 Misc : SFAM SURROGATE
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP6631

Quant Time: Sep 17 04:17:30 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG090924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 10 01:28:17 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	30364	20.000	ng/u1	0.00
20) Naphthalene-d8	10.691	136	115568	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	104312	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.271	188	302944	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.519	240	360524	20.000	ng/u1	# 0.00
88) Perylene-d12	24.586	264	409544	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	9154	13.628	ng/uL	0.00
4) Pyridine-d5	3.764	84	133677	62.519	ng/u1	0.00
7) Phenol-d5	7.072	99	157055	58.986	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.230	67	85808	53.754	ng/u1	0.00
11) 2-Chlorophenol-d4	7.436	132	125825	64.878	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	137956	62.176	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	67233	81.398	ng/u1	0.00
24) 2-Nitrophenol-d4	9.774	143	78384	86.772	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.315	165	154315	78.442	ng/u1	0.00
31) 4-Chloroaniline-d4	10.826	131	189963	70.249	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	573280	71.368	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	625434	70.106	ng/u1	0.00
54) 4-Nitrophenol-d4	14.733	143	92620	67.874	ng/u1	0.00
60) Fluorene-d10	15.521	176	522140	72.134	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.638	200	127425	73.124	ng/u1	0.00
73) Anthracene-d10	17.371	188	1011877	70.775	ng/u1	0.00
81) Pyrene-d10	19.669	212	1337300	65.917	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.387	264	1560656	72.169	ng/u1	0.00

Target Compounds	Qvalue

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

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