

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111524\
 Data File : BG063447.D
 Acq On : 16 Nov 2024 5:03
 Operator : RC/JU
 Sample : P4803-14
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A0AQ9

Quant Time: Nov 16 04:38:31 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 15:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.830	152	86755	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.620	136	379059	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.469	164	309875	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.224	188	740767	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.478	240	783557	20.000	ng/u1	# 0.00
88) Perylene-d12	24.516	264	840141	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.300	96	15770	8.219	ng/uL	0.00
4) Pyridine-d5	3.711	84	103747	16.360	ng/u1	0.00
7) Phenol-d5	7.007	99	148855	19.015	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.166	67	182906	39.833	ng/u1	0.00
11) 2-Chlorophenol-d4	7.365	132	196328	38.638	ng/u1	-0.01
15) 4-Methylphenol-d8	8.546	113	230767	34.779	ng/u1	0.00
21) Nitrobenzene-d5	8.987	128	113259	42.500	ng/u1	0.00
24) 2-Nitrophenol-d4	9.704	143	139139	40.398	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.250	165	269403	39.857	ng/u1	0.00
31) 4-Chloroaniline-d4	10.756	131	246761	28.029	ng/u1	0.00
46) Dimethylphthalate-d6	13.875	166	927712	38.221	ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	1002643	42.203	ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	50179	15.099	ng/u1	0.01
60) Fluorene-d10	15.468	176	823936	40.221	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.591	200	183082	39.326	ng/u1	0.00
73) Anthracene-d10	17.324	188	1449651	45.533	ng/u1	0.00
81) Pyrene-d10	19.622	212	1880007	47.855	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.310	264	1898000	46.791	ng/u1	0.00

Target Compounds Qvalue

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

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