

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032025\
 Data File : BG064117.D
 Acq On : 20 Mar 2025 14:36
 Operator : RC/JU
 Sample : PB167183BL
 Misc : MDL-WATER 4 ppm
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK183

Quant Time: Mar 20 15:51:12 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031425.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 14 14:37:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.854	152	28486	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.644	136	132535	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.481	164	97319	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.225	188	232862	20.000	ng/u1	0.00
79) Chrysene-d12	21.455	240	228411	20.000	ng/u1	0.00
88) Perylene-d12	24.475	264	241029	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.335	96	4350	6.114	ng/uL	0.00
4) Pyridine-d5	3.735	84	68645	33.581	ng/u1	0.00
7) Phenol-d5	7.019	99	94142	36.525	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.190	67	55109	33.931	ng/u1	0.00
11) 2-Chlorophenol-d4	7.389	132	72110	35.058	ng/u1	0.00
15) 4-Methylphenol-d8	8.553	113	76314	34.904	ng/u1	0.00
21) Nitrobenzene-d5	9.005	128	32665	32.861	ng/u1	0.00
24) 2-Nitrophenol-d4	9.728	143	34248	32.806	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.262	165	71049	33.751	ng/u1	0.00
31) 4-Chloroaniline-d4	10.774	131	100853	34.168	ng/u1	0.00
46) Dimethylphthalate-d6	13.894	166	267888	35.074	ng/u1	0.00
49) Acenaphthylene-d8	14.176	160	289472	34.081	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	39661	30.905	ng/u1	0.00
60) Fluorene-d10	15.474	176	232475	35.311	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	30301	24.129	ng/u1	0.00
73) Anthracene-d10	17.325	188	394532	33.787	ng/u1	0.00
81) Pyrene-d10	19.610	212	449253	35.651	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.270	264	458518	36.501	ng/u1	0.00

Target Compounds Qvalue

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

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