

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062845.D
 Acq On : 15 Sep 2024 14:18
 Operator : JU/RC
 Sample : PB163398BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK398

Quant Time: Sep 15 23:34:20 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	34680	20.000	ng/u1	0.00
20) Naphthalene-d8	10.691	136	143071	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.528	164	122532	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.272	188	343474	20.000	ng/u1	0.00
79) Chrysene-d12	21.520	240	388804	20.000	ng/u1	0.00
88) Perylene-d12	24.581	264	436115	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.364	96	5317	6.931	ng/uL	0.00
4) Pyridine-d5	3.764	84	77070	31.559	ng/u1	0.00
7) Phenol-d5	7.066	99	89449	29.414	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.230	67	53455	29.319	ng/u1	0.00
11) 2-Chlorophenol-d4	7.436	132	71788	32.409	ng/u1	0.00
15) 4-Methylphenol-d8	8.605	113	80166	31.634	ng/u1	0.00
21) Nitrobenzene-d5	9.052	128	37888	37.053	ng/u1	0.00
24) 2-Nitrophenol-d4	9.774	143	44396	39.699	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.315	165	85747	35.208	ng/u1	0.00
31) 4-Chloroaniline-d4	10.826	131	107507	32.114	ng/u1	0.00
46) Dimethylphthalate-d6	13.934	166	318890	33.796	ng/u1	0.00
49) Acenaphthylene-d8	14.222	160	350724	33.467	ng/u1	0.00
54) 4-Nitrophenol-d4	14.728	143	49711	31.013	ng/u1	0.00
60) Fluorene-d10	15.521	176	295403	34.742	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.638	200	61759	31.259	ng/u1	0.00
73) Anthracene-d10	17.371	188	573715	35.393	ng/u1	0.00
81) Pyrene-d10	19.663	212	755005	34.508	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.375	264	797878	34.648	ng/u1	0.00

Target Compounds Qvalue

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

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