

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091324\
 Data File : BG062782.D
 Acq On : 13 Sep 2024 11:07
 Operator : JU/RC
 Sample : PB163356BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK356

Quant Time: Sep 13 23:07:02 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	7.899	152	45173	20.000	ng/u1	0.00
20) Naphthalene-d8	10.690	136	189670	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.527	164	161883	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.270	188	429487	20.000	ng/u1	0.00
79) Chrysene-d12	21.518	240	487104	20.000	ng/u1	0.00
88) Perylene-d12	24.574	264	535084	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.357	96	7699	7.704	ng/uL	0.00
4) Pyridine-d5	3.763	84	113037	35.535	ng/u1	0.00
7) Phenol-d5	7.065	99	113703	28.705	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.229	67	68659	28.911	ng/u1	0.00
11) 2-Chlorophenol-d4	7.429	132	87350	30.274	ng/u1	0.00
15) 4-Methylphenol-d8	8.598	113	97719	29.603	ng/u1	-0.01
21) Nitrobenzene-d5	9.051	128	45403	33.493	ng/u1	0.00
24) 2-Nitrophenol-d4	9.773	143	49759	33.563	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.314	165	100005	30.974	ng/u1	0.00
31) 4-Chloroaniline-d4	10.825	131	130283	29.356	ng/u1	0.00
46) Dimethylphthalate-d6	13.933	166	402284	32.270	ng/u1	0.00
49) Acenaphthylene-d8	14.221	160	434428	31.378	ng/u1	0.00
54) 4-Nitrophenol-d4	14.726	143	63685	30.073	ng/u1	0.00
60) Fluorene-d10	15.520	176	380023	33.829	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.637	200	67144	27.178	ng/u1	0.00
73) Anthracene-d10	17.370	188	718759	35.461	ng/u1	0.00
81) Pyrene-d10	19.662	212	969634	35.374	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.368	264	968435	34.276	ng/u1	-0.01

Target Compounds	Qvalue

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

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