

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062224\
 Data File : BG061651.D
 Acq On : 22 Jun 2024 16:28
 Operator : MA/JU
 Sample : PB161462BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK462

Quant Time: Jun 23 00:19:56 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 21 05:03:55 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.079	152	62510	20.000	ng/u1	0.00
20) Naphthalene-d8	10.888	136	314857	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.701	164	220779	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.445	188	505123	20.000	ng/u1	0.00
79) Chrysene-d12	21.705	240	436590	20.000	ng/u1	0.00
88) Perylene-d12	24.930	264	535757	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.467	96	10600	6.126	ng/uL	0.00
4) Pyridine-d5	3.884	84	154435	29.518	ng/u1	0.00
7) Phenol-d5	7.216	99	208641	30.745	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.398	67	124344	29.394	ng/u1	0.00
11) 2-Chlorophenol-d4	7.598	132	144392	31.007	ng/u1	0.00
15) 4-Methylphenol-d8	8.767	113	161007	30.691	ng/u1	0.00
21) Nitrobenzene-d5	9.243	128	75035	35.098	ng/u1	0.00
24) 2-Nitrophenol-d4	9.960	143	83767	38.687	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.494	165	160995	31.401	ng/u1	0.00
31) 4-Chloroaniline-d4	11.017	131	227049	29.250	ng/u1	0.00
46) Dimethylphthalate-d6	14.108	166	528853	31.128	ng/u1	0.00
49) Acenaphthylene-d8	14.396	160	612837	32.914	ng/u1	0.00
54) 4-Nitrophenol-d4	14.872	143	86515	29.650	ng/u1	0.00
60) Fluorene-d10	15.694	176	467927	31.460	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.794	200	59668	27.625	ng/u1	0.00
73) Anthracene-d10	17.545	188	732612	31.774	ng/u1	0.00
81) Pyrene-d10	19.825	212	802875	32.922	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.713	264	821386	32.100	ng/u1	0.00

Target Compounds Qvalue

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

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