

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081924\
 Data File : BG062461.D
 Acq On : 19 Aug 2024 12:49
 Operator : MA/JU
 Sample : PB162773BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK773

Quant Time: Aug 20 00:24:09 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG081424.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 14 15:31:21 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.947	152	59226	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.737	136	287632	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.567	164	232682	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.305	188	564434	20.000	ng/u1	-0.01
79) Chrysene-d12	21.546	240	536839	20.000	ng/u1	-0.02
88) Perylene-d12	24.630	264	610330	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.406	96	7976	5.113	ng/uL	-0.01
4) Pyridine-d5	3.811	84	113889	23.959	ng/u1	-0.01
7) Phenol-d5	7.101	99	146844	24.923	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.271	67	94889	25.567	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.477	132	93873	23.324	ng/u1	-0.01
15) 4-Methylphenol-d8	8.640	113	107498	21.766	ng/u1	-0.02
21) Nitrobenzene-d5	9.098	128	48439	26.260	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.821	143	50406	27.470	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.355	165	106020	22.685	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.866	131	151271	20.410	ng/u1	-0.01
46) Dimethylphthalate-d6	13.974	166	433067	23.927	ng/u1	-0.01
49) Acenaphthylene-d8	14.256	160	451594	22.535	ng/u1	-0.02
54) 4-Nitrophenol-d4	14.749	143	64542	22.354	ng/u1	-0.01
60) Fluorene-d10	15.554	176	380903	23.334	ng/u1	-0.02
65) 4,6-Dinitro-2-methylph...	15.666	200	55420	21.610	ng/u1	-0.02
73) Anthracene-d10	17.405	188	639487	23.585	ng/u1	-0.01
81) Pyrene-d10	19.690	212	817298	25.996	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.424	264	739890	22.843	ng/u1	-0.02

Target Compounds Qvalue

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

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