

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081424\
 Data File : BG062400.D
 Acq On : 14 Aug 2024 18:20
 Operator : MA/JU
 Sample : P3589-03
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E0CK3

Quant Time: Aug 15 00:54:56 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	7.959	152	48898	20.000	ng/u1	0.00
20) Naphthalene-d8	10.750	136	227003	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.574	164	187035	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.317	188	475491	20.000	ng/u1	0.00
79) Chrysene-d12	21.558	240	491065	20.000	ng/u1	0.00
88) Perylene-d12	24.648	264	565932	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.418	96	4557	3.538	ng/uL	0.00
4) Pyridine-d5	3.830	84	21262	5.418	ng/u1	0.00
7) Phenol-d5	7.113	99	17437	3.585	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.284	67	32465	10.595	ng/u1	0.00
11) 2-Chlorophenol-d4	7.489	132	30556	9.196	ng/u1	0.00
15) 4-Methylphenol-d8	8.652	113	31727	7.781	ng/u1	0.00
21) Nitrobenzene-d5	9.105	128	17497	12.019	ng/u1	0.00
24) 2-Nitrophenol-d4	9.827	143	13328	9.203	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.368	165	35185	9.539	ng/u1	0.00
31) 4-Chloroaniline-d4	10.879	131	42218	7.217	ng/u1	0.00
46) Dimethylphthalate-d6	13.980	166	172834	11.880	ng/u1	0.00
49) Acenaphthylene-d8	14.268	160	169291	10.509	ng/u1	0.00
54) 4-Nitrophenol-d4	14.762	143	6010	2.590	ng/u1	0.00
60) Fluorene-d10	15.567	176	148913	11.349	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.672	200	10304	4.769	ng/u1	-0.01
73) Anthracene-d10	17.411	188	269121	11.782	ng/u1	0.00
81) Pyrene-d10	19.696	212	355087	12.347	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.431	264	351562	11.706	ng/u1	-0.01

Target Compounds Qvalue

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

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