

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG073124\
 Data File : BG062224.D
 Acq On : 31 Jul 2024 19:04
 Operator : MA/JU
 Sample : P3398-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 E08K6

Quant Time: Aug 01 04:05:18 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG073124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 01 03:46:07 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.979	152	34305	20.000	ng/u1	0.00
20) Naphthalene-d8	10.775	136	159565	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.599	164	120096	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.337	188	302213	20.000	ng/u1	0.00
79) Chrysene-d12	21.584	240	333813	20.000	ng/u1	0.00
88) Perylene-d12	24.697	264	374564	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.438	96	2950	3.660	ng/uL	0.00
4) Pyridine-d5	3.855	84	4924	2.000	ng/u1	0.01
7) Phenol-d5	7.127	99	21272	6.479	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.309	67	57997	28.645	ng/u1	0.00
11) 2-Chlorophenol-d4	7.509	132	58357	26.427	ng/u1	0.00
15) 4-Methylphenol-d8	8.666	113	45305	16.580	ng/u1	0.00
21) Nitrobenzene-d5	9.130	128	27511	41.569	ng/u1	0.00
24) 2-Nitrophenol-d4	9.853	143	26428	42.624	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	69486	29.930	ng/u1	0.00
31) 4-Chloroaniline-d4	10.898	131	52279	13.309	ng/u1	0.00
46) Dimethylphthalate-d6	14.006	166	302445	33.417	ng/u1	0.00
49) Acenaphthylene-d8	14.288	160	309897	31.016	ng/u1	0.00
54) 4-Nitrophenol-d4	14.764	143	7632	7.026	ng/u1	-0.01
60) Fluorene-d10	15.592	176	258032	32.086	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	20002	27.921	ng/u1	0.00
73) Anthracene-d10	17.436	188	467806	33.590	ng/u1	0.00
81) Pyrene-d10	19.722	212	628729	33.654	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.486	264	662400	36.174	ng/u1	0.00

Target Compounds	Qvalue

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

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