

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG062022\
 Data File : BG053960.D
 Acq On : 20 Jun 2022 22:33
 Operator : CG/JU
 Sample : N3358-17
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 EW4T0

Quant Time: Jun 20 23:59:08 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG062022.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jun 20 15:29:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.079	152	18817	20.000	ng/u1	0.00
20) Naphthalene-d8	10.887	136	76922	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.700	164	59938	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.444	188	165518	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.716	240	198945	20.000	ng/u1	# 0.00
88) Perylene-d12	24.982	264	202791	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.502	96	1762	4.455	ng/uL	0.00
4) Pyridine-d5	3.919	84	16853	15.667	ng/u1	0.00
7) Phenol-d5	7.238	99	926	0.687	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.397	67	26445	32.520	ng/u1	0.00
11) 2-Chlorophenol-d4	0.000	132	0	0.000	ng/u1	
15) 4-Methylphenol-d8	8.778	113	6229	5.419	ng/u1	0.00
21) Nitrobenzene-d5	9.236	128	18447	32.309	ng/u1	0.00
24) 2-Nitrophenol-d4	0.000	143	0	0.000	ng/u1	
28) 2,4-Dichlorophenol-d3	0.000	165	0	0.000	ng/u1	
31) 4-Chloroaniline-d4	11.016	131	50765	29.939	ng/u1	0.00
46) Dimethylphthalate-d6	14.101	166	157730	33.451	ng/u1	0.00
49) Acenaphthylene-d8	14.395	160	167278	32.625	ng/u1	0.00
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/u1	
60) Fluorene-d10	15.693	176	142828	33.722	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0	0.000	ng/u1	
73) Anthracene-d10	17.544	188	254501	34.047	ng/u1	0.00
81) Pyrene-d10	19.830	212	340853	32.860	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.753	264	353559	34.540	ng/u1	0.00

Target Compounds	Qvalue

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

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