

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110423\
 Data File : BG059624.D
 Acq On : 4 Nov 2023 15:58
 Operator : MA/JU
 Sample : 04996-11
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A49K6

Quant Time: Nov 05 23:54:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 04 02:46:50 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.356	152	38362	20.000	ng/u1	0.00
20) Naphthalene-d8	11.199	136	190970	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.983	164	133683	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.733	188	232979	20.000	ng/u1	0.00
79) Chrysene-d12	22.063	240	263238	20.000	ng/u1	# 0.00
88) Perylene-d12	25.671	264	302400	20.000	ng/u1	# 0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.649	96	2886	2.820	ng/uL	0.00
4) Pyridine-d5	4.078	84	22202	7.376	ng/u1	0.00
7) Phenol-d5	7.468	99	21987	5.606	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.668	67	51181	24.407	ng/u1	0.00
11) 2-Chlorophenol-d4	7.868	132	58364	21.911	ng/u1	0.00
15) 4-Methylphenol-d8	9.037	113	42070	13.106	ng/u1	0.00
21) Nitrobenzene-d5	9.548	128	39683	28.414	ng/u1	0.00
24) 2-Nitrophenol-d4	10.265	143	46379	27.518	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.800	165	77388	25.844	ng/u1	0.00
31) 4-Chloroaniline-d4	11.340	131	127429	27.043	ng/u1	0.00
46) Dimethylphthalate-d6	14.372	166	379680	35.389	ng/u1	0.00
49) Acenaphthylene-d8	14.683	160	418508	32.713	ng/u1	0.00
54) 4-Nitrophenol-d4	15.148	143	12160	6.419	ng/u1	0.00
60) Fluorene-d10	15.970	176	339779	33.639	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.070	200	64734	43.898	ng/u1	0.00
73) Anthracene-d10	17.833	188	407193	33.394	ng/u1	0.00
81) Pyrene-d10	20.107	212	540235	32.070	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.418	264	592437	33.466	ng/u1	0.00

Target Compounds	Qvalue

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

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