

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG103123\
 Data File : BG059566.D
 Acq On : 31 Oct 2023 18:31
 Operator : MA/JU
 Sample : 04950-16
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 A49E8

Quant Time: Nov 01 03:20:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG102723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 28 03:56:42 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.352	152	77441	20.000	ng/ul	# 0.00
20) Naphthalene-d8	11.202	136	413138	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.986	164	260784	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.735	188	559891	20.000	ng/ul	0.00
79) Chrysene-d12	22.072	240	466009	20.000	ng/ul	# 0.00
88) Perylene-d12	25.679	264	501819	20.000	ng/ul	#-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.652	96	5588	2.700	ng/uL	0.00
4) Pyridine-d5	4.093	84	15329	2.650	ng/ul	0.00
7) Phenol-d5	7.465	99	38265	4.808	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.677	67	84888	23.224	ng/ul	0.00
11) 2-Chlorophenol-d4	7.871	132	111212	17.226	ng/ul	0.00
15) 4-Methylphenol-d8	9.034	113	73429	11.630	ng/ul	0.00
21) Nitrobenzene-d5	9.551	128	76808	26.147	ng/ul	0.00
24) 2-Nitrophenol-d4	10.274	143	93980	31.164	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.802	165	160438	24.147	ng/ul	0.00
31) 4-Chloroaniline-d4	11.343	131	34250	3.131	ng/ul	0.00
46) Dimethylphthalate-d6	14.381	166	630175	26.673	ng/ul	0.00
49) Acenaphthylene-d8	14.686	160	719044	25.933	ng/ul	0.00
54) 4-Nitrophenol-d4	15.150	143	22656	5.592	ng/ul	0.00
60) Fluorene-d10	15.973	176	534225	26.458	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.073	200	81721	30.821	ng/ul	0.00
73) Anthracene-d10	17.835	188	893347	29.451	ng/ul	0.00
81) Pyrene-d10	20.109	212	955483	29.498	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.426	264	914968	30.964	ng/ul	-0.01

Target Compounds Qvalue

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

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