

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063641.D
 Acq On : 10 Dec 2024 21:05
 Operator : RC/JU
 Sample : P5109-16
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 10 23:27:19 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG120624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 06 15:14:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.812	152	122401	20.000	ng/ul	0.00
20) Naphthalene-d8	10.603	136	558466	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.452	164	410511	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.201	188	957997	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.455	240	886115	20.000	ng/ul	#-0.02
88) Perylene-d12	24.475	264	986296	20.000	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.282	96	17317	6.246	ng/uL	0.00
4) Pyridine-d5	3.694	84	62384	6.082	ng/ul	0.00
7) Phenol-d5	6.996	99	101204	8.398	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.143	67	308031	37.803	ng/ul	0.00
11) 2-Chlorophenol-d4	7.348	132	253621	35.726	ng/ul	0.00
15) 4-Methylphenol-d8	8.523	113	192571	21.552	ng/ul	0.00
21) Nitrobenzene-d5	9.111	128	247	0.063	ng/ul	0.14
24) 2-Nitrophenol-d4	9.687	143	170033	37.905	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.233	165	321680	33.305	ng/ul	0.00
31) 4-Chloroaniline-d4	10.738	131	331097	28.366	ng/ul	0.00
46) Dimethylphthalate-d6	13.858	166	1187813	39.813	ng/ul	0.00
49) Acenaphthylene-d8	14.146	160	1195742	36.666	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	16182	4.236	ng/ul	0.00
60) Fluorene-d10	15.451	176	1013582	34.465	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.580	200	144686	26.735	ng/ul	0.00
73) Anthracene-d10	17.301	188	1910140	43.667	ng/ul	-0.01
81) Pyrene-d10	19.604	212	2268893	46.194	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.275	264	2246048	45.341	ng/ul	-0.02

Target Compounds	Qvalue

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

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