

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121024\
 Data File : BG063637.D
 Acq On : 10 Dec 2024 18:34
 Operator : RC/JU
 Sample : P5109-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 10 23:26:41 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	7.811	152	149181	20.000	ng/ul	0.00
20) Naphthalene-d8	10.601	136	641583	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.450	164	478165	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.206	188	1095687	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.459	240	971640	20.000	ng/ul	#-0.01
88) Perylene-d12	24.479	264	1060377	20.000	ng/ul	-0.03

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	19190	5.679	ng/uL	0.00
4) Pyridine-d5	3.692	84	81215	6.497	ng/ul	0.00
7) Phenol-d5	6.994	99	134673	9.169	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.147	67	354338	35.680	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	300873	34.774	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	244229	22.427	ng/ul	0.00
21) Nitrobenzene-d5	8.968	128	174796	38.814	ng/ul	0.00
24) 2-Nitrophenol-d4	9.685	143	195377	37.912	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.231	165	372111	33.536	ng/ul	0.00
31) 4-Chloroaniline-d4	10.737	131	50226	3.746	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	1356092	39.022	ng/ul	0.00
49) Acenaphthylene-d8	14.144	160	1359733	35.796	ng/ul	0.00
54) 4-Nitrophenol-d4	14.679	143	22171	4.982	ng/ul	0.00
60) Fluorene-d10	15.449	176	1172410	34.225	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.578	200	161838	26.146	ng/ul	0.00
73) Anthracene-d10	17.300	188	2005955	40.095	ng/ul	-0.01
81) Pyrene-d10	19.603	212	2255459	41.878	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.274	264	2226484	41.806	ng/ul	-0.02

Target Compounds Qvalue

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

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