

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012924\
 Data File : BG060472.D
 Acq On : 29 Jan 2024 21:41
 Operator : MA/JU
 Sample : P1239-17
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0FC7

Quant Time: Jan 30 01:49:23 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG012624.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jan 26 06:06:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	168961	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	845740	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.564	164	581723	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.308	188	1498125	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	1475359	20.000	ng/ul	0.00
88) Perylene-d12	24.735	264	1622788	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.366	96	15322	4.544	ng/uL	0.00
4) Pyridine-d5	3.777	84	105641	9.904	ng/ul	0.00
7) Phenol-d5	7.091	99	80750	5.032	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.249	67	253218	25.112	ng/ul	0.00
11) 2-Chlorophenol-d4	7.455	132	214784	20.044	ng/ul	0.00
15) 4-Methylphenol-d8	8.630	113	173600	13.134	ng/ul	0.00
21) Nitrobenzene-d5	9.153	128	1263	0.202	ng/ul	0.06
24) 2-Nitrophenol-d4	9.805	143	184308	25.286	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	353966	23.350	ng/ul	0.00
31) 4-Chloroaniline-d4	10.863	131	438069	22.029	ng/ul	0.00
46) Dimethylphthalate-d6	13.965	166	1448921	30.439	ng/ul	0.00
49) Acenaphthylene-d8	14.253	160	1552071	29.068	ng/ul	0.00
54) 4-Nitrophenol-d4	14.782	143	11571	1.697	ng/ul	0.02
60) Fluorene-d10	15.557	176	1340258	32.133	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	20595	2.220	ng/ul	0.00
73) Anthracene-d10	17.408	188	2355680	33.171	ng/ul	0.00
81) Pyrene-d10	19.700	212	2923799	33.065	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.517	264	2942544	34.566	ng/ul	0.00

Target Compounds Qvalue

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

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