

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023093.D
 Acq On : 13 Nov 2024 22:56
 Operator : RC/JU
 Sample : P4724-07
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Nov 14 01:37:25 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 13 04:49:01 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	85293	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	335520	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.192	164	266264	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.004	188	665034	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.445	240	703913	20.000	ng/ul	0.00
88) Perylene-d12	24.657	264	825094	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	5385	3.035	ng/uL	0.00
4) Pyridine-d5	3.558	84	61259	11.620	ng/ul	0.01
7) Phenol-d5	6.787	99	105195	16.682	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	63083	17.027	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	77297	16.737	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	86928	16.752	ng/ul	0.00
21) Nitrobenzene-d5	8.728	128	40666	17.410	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	48197	17.308	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	99746	17.104	ng/ul	0.01
31) 4-Chloroaniline-d4	10.487	131	115041	16.536	ng/ul	0.00
46) Dimethylphthalate-d6	13.598	166	380289	19.817	ng/ul	-0.01
49) Acenaphthylene-d8	13.881	160	399042	19.928	ng/ul	-0.02
54) 4-Nitrophenol-d4	14.481	143	38036	13.189	ng/ul	0.03
60) Fluorene-d10	15.216	176	342480	20.355	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.357	200	46993	11.796	ng/ul	0.00
73) Anthracene-d10	17.104	188	580997	20.633	ng/ul	-0.01
81) Pyrene-d10	19.486	212	725790	22.385	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.445	264	823188	20.945	ng/ul	0.00

Target Compounds Qvalue

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

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