

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012229.D
 Acq On : 20 Oct 2022 23:57
 Operator : CG/JU
 Sample : N5103-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBB0

Quant Time: Oct 21 01:36:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.822	152	451403	20.000 ng/u1	0.00
20) Naphthalene-d8	10.628	136	1946236	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1199695	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.234	188	2288446	20.000 ng/u1	0.00
79) Chrysene-d12	21.327	240	1633560	20.000 ng/u1	0.00
88) Perylene-d12	23.727	264	1667863	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.258	96	41139	3.688 ng/uL	0.00
4) Pyridine-d5	3.681	84	268292	8.347 ng/u1	0.00
7) Phenol-d5	6.993	99	213105	5.468 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	673885	28.513 ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	651853	21.975 ng/u1	0.00
15) 4-Methylphenol-d8	8.540	113	415511	13.719 ng/u1	0.00
21) Nitrobenzene-d5	8.993	128	456938	31.718 ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	469569	30.211 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.252	165	764263	26.214 ng/u1	0.00
31) 4-Chloroaniline-d4	10.775	131	880377	21.253 ng/u1	0.00
46) Dimethylphthalate-d6	13.887	166	2664858	31.933 ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	2960982	31.084 ng/u1	0.00
54) 4-Nitrophenol-d4	14.693	143	55995	3.582 ng/u1	0.00
60) Fluorene-d10	15.469	176	2271344	31.637 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	287076	22.051 ng/u1	0.00
73) Anthracene-d10	17.334	188	3588567	35.775 ng/u1	0.00
81) Pyrene-d10	19.575	212	3533151	37.988 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	3049610	36.461 ng/u1	-0.01

Target Compounds	Qvalue

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

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