

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012237.D
 Acq On : 21 Oct 2022 10:21
 Operator : CG/JU
 Sample : N5103-16
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BA1

Quant Time: Oct 21 22:29:47 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	345191	20.000 ng/u1	0.00
20) Naphthalene-d8	10.628	136	1485236	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.469	164	895825	20.000 ng/u1	-0.01
64) Phenanthrene-d10	17.233	188	1701768	20.000 ng/u1	0.00
79) Chrysene-d12	21.327	240	1261074	20.000 ng/u1	0.00
88) Perylene-d12	23.727	264	1329504	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.258	96	34452	4.039 ng/uL	0.00
4) Pyridine-d5	3.681	84	342259	13.925 ng/u1	0.00
7) Phenol-d5	6.993	99	446284	14.976 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	725281	40.130 ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	675303	29.770 ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	561806	24.256 ng/u1	-0.01
21) Nitrobenzene-d5	8.993	128	490793	44.643 ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	324894	27.391 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	697229	31.337 ng/u1	-0.01
31) 4-Chloroaniline-d4	10.769	131	1145273	36.230 ng/u1	-0.01
46) Dimethylphthalate-d6	13.887	166	2843389	45.629 ng/u1	0.00
49) Acenaphthylene-d8	14.169	160	3254282	45.752 ng/u1	0.00
54) 4-Nitrophenol-d4	14.692	143	55036	4.714 ng/u1	0.00
60) Fluorene-d10	15.469	176	2448698	45.676 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	24125	2.492 ng/u1	0.00
73) Anthracene-d10	17.334	188	3766080	50.488 ng/u1	0.00
81) Pyrene-d10	19.575	212	3701293	51.550 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.574	264	3488665	52.325 ng/u1	-0.01

Target Compounds Qvalue

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

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