

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012230.D
 Acq On : 21 Oct 2022 00:31
 Operator : CG/JU
 Sample : N5103-06
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BX3

Quant Time: Oct 21 01:38:26 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	398999	20.000 ng/u1	0.00
20) Naphthalene-d8	10.628	136	1713783	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.475	164	1003043	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1803073	20.000 ng/u1	0.00
79) Chrysene-d12	21.321	240	1437365	20.000 ng/u1	-0.01
88) Perylene-d12	23.727	264	1549681	20.000 ng/u1	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.258	96	45556	4.621 ng/uL	0.00
4) Pyridine-d5	3.687	84	172037	6.056 ng/u1	0.00
7) Phenol-d5	6.993	99	220888	6.413 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	664796	31.823 ng/u1	0.00
11) 2-Chlorophenol-d4	7.352	132	655730	25.009 ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	406459	15.182 ng/u1	-0.01
21) Nitrobenzene-d5	8.993	128	443413	34.954 ng/u1	0.00
24) 2-Nitrophenol-d4	9.710	143	458297	33.485 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	728914	28.393 ng/u1	0.00
31) 4-Chloroaniline-d4	10.769	131	716701	19.649 ng/u1	-0.01
46) Dimethylphthalate-d6	13.887	166	2372919	34.009 ng/u1	0.00
49) Acenaphthylene-d8	14.163	160	2706996	33.989 ng/u1	-0.01
54) 4-Nitrophenol-d4	14.698	143	51381	3.931 ng/u1	0.00
60) Fluorene-d10	15.469	176	2000546	33.328 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	251971	24.565 ng/u1	0.00
73) Anthracene-d10	17.333	188	3005523	38.028 ng/u1	0.00
81) Pyrene-d10	19.569	212	3012506	36.811 ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.574	264	3019218	38.850 ng/u1	-0.01

Target Compounds Qvalue

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

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