

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012064.D
 Acq On : 12 Oct 2022 05:26
 Operator : CG/JU
 Sample : N4991-10
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BG2

Quant Time: Oct 12 07:03:56 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 23:54:43 2022
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.852	152	251171	20.000 ng/u1	0.00
20) Naphthalene-d8	10.663	136	1006437	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.504	164	564460	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.257	188	1173398	20.000 ng/u1	-0.01
79) Chrysene-d12	21.345	240	1276530	20.000 ng/u1	# 0.00
88) Perylene-d12	23.763	264	1385337	20.000 ng/u1	0.00

System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.275	96	28377	4.454 ng/uL	0.00
4) Pyridine-d5	3.705	84	225432	13.306 ng/u1	0.00
7) Phenol-d5	7.022	99	358691	17.756 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	282819	24.665 ng/u1	0.00
11) 2-Chlorophenol-d4	7.381	132	321477	20.049 ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	257336	16.585 ng/u1	0.00
21) Nitrobenzene-d5	9.022	128	196502	27.645 ng/u1	0.00
24) 2-Nitrophenol-d4	9.746	143	167912	23.297 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.281	165	281532	20.155 ng/u1	0.00
31) 4-Chloroaniline-d4	10.804	131	288034	13.993 ng/u1	0.00
46) Dimethylphthalate-d6	13.916	166	1183159	33.097 ng/u1	0.00
49) Acenaphthylene-d8	14.192	160	1511437	34.191 ng/u1	0.00
54) 4-Nitrophenol-d4	14.710	143	88508	12.611 ng/u1	0.00
60) Fluorene-d10	15.498	176	1262435	39.032 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	85852	13.856 ng/u1	0.00
73) Anthracene-d10	17.357	188	1728069	33.915 ng/u1	0.00
81) Pyrene-d10	19.598	212	2047898	31.315 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.610	264	1388310	20.615 ng/u1	0.00

Target Compounds	Qvalue

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

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