

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012112.D
 Acq On : 13 Oct 2022 10:07
 Operator : CG/JU
 Sample : N5026-10
 Misc :
 ALS Vial : 83 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :

Quant Time: Oct 13 23:30:48 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	227759	20.000	ng/ul	0.00
20) Naphthalene-d8	10.651	136	925968	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.498	164	511083	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	1055548	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1113172	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264	1176741	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	26066	4.512	ng/uL	0.00
4) Pyridine-d5	3.705	84	110975	7.224	ng/ul	0.00
7) Phenol-d5	7.016	99	162018	8.845	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	393292	37.826	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	465693	32.028	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	285946	20.324	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	274579	41.987	ng/ul	0.00
24) 2-Nitrophenol-d4	9.740	143	284857	42.958	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	468238	36.435	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	475731	25.120	ng/ul	0.00
46) Dimethylphthalate-d6	13.910	166	1384900	42.786	ng/ul	0.00
49) Acenaphthylene-d8	14.186	160	1661298	41.506	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	39624	6.236	ng/ul	0.00
60) Fluorene-d10	15.492	176	1268550	43.318	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.616	200	184470	33.097	ng/ul	0.00
73) Anthracene-d10	17.357	188	2024067	44.159	ng/ul	0.00
81) Pyrene-d10	19.592	212	2476042	43.418	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.604	264	2608404	45.598	ng/ul	0.00

Target Compounds Qvalue

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

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