

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018110.D
 Acq On : 12 Nov 2023 17:45
 Operator : MA/JU
 Sample : 05082-05
 Misc :
 ALS Vial : 90 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH331

Quant Time: Nov 12 22:38:42 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	78404	20.000	ng/u1	0.00
20) Naphthalene-d8	10.663	136	309395	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.516	164	180284	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.298	188	406426	20.000	ng/u1	-0.01
79) Chrysene-d12	21.427	240	417443	20.000	ng/u1	0.00
88) Perylene-d12	23.927	264	480474	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	9225	4.151	ng/uL	0.00
4) Pyridine-d5	3.669	84	43431	7.494	ng/u1	0.00
7) Phenol-d5	7.016	99	68255	9.017	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	136140	30.605	ng/u1	0.00
11) 2-Chlorophenol-d4	7.369	132	145497	24.273	ng/u1	0.00
15) 4-Methylphenol-d8	8.569	113	110620	17.834	ng/u1	0.00
21) Nitrobenzene-d5	9.051	128	85790	30.777	ng/u1	0.00
24) 2-Nitrophenol-d4	9.763	143	93313	29.549	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	160734	28.557	ng/u1	0.00
31) 4-Chloroaniline-d4	10.845	131	192014	24.877	ng/u1	0.00
46) Dimethylphthalate-d6	13.939	166	593365	35.671	ng/u1	-0.01
49) Acenaphthylene-d8	14.216	160	607959	33.556	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.828	143	15624	6.286	ng/u1	0.05
60) Fluorene-d10	15.522	176	491773	35.808	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	148	0.050	ng/u1	0.11
73) Anthracene-d10	17.398	188	857669	39.018	ng/u1	-0.01
81) Pyrene-d10	19.657	212	1073135	38.486	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.768	264	1148115	41.095	ng/u1	0.00

Target Compounds	Qvalue

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

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